

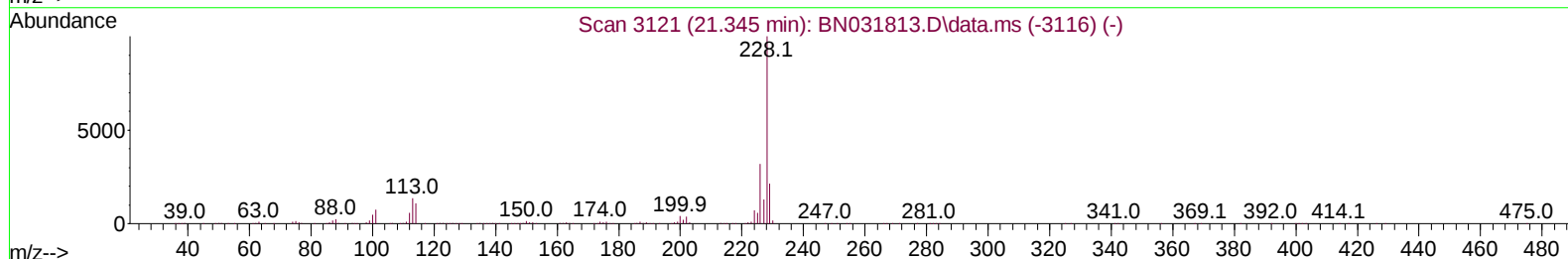
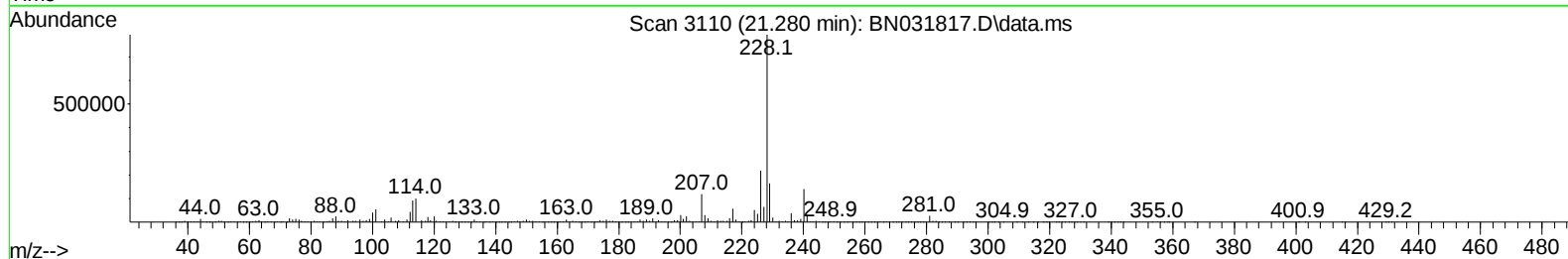
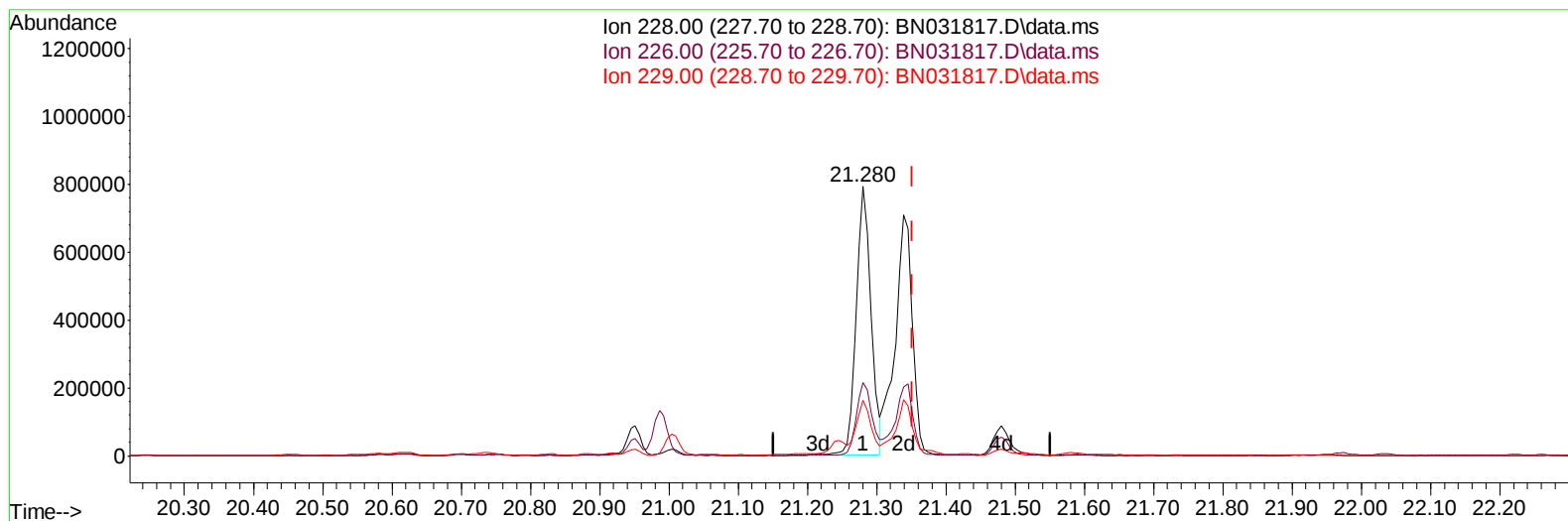
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060624\
Data File : BN031817.D
Acq On : 06 Jun 2024 13:55
Operator : MA/JU
Sample : P2634-20
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHB4

Manual IntegrationsAPPROVED

Quant Time: Jun 06 14:34:57 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 29 15:13:07 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/07/2024
Supervised By :mohammad ahmed 06/08/2024



TIC: BN031817.D\data.ms

(87) Chrysene

21.280min (-0.071) 12.87 ng/u1

response 1162349

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	29.60	27.32
229.00	18.90	20.68
0.00	0.00	0.00

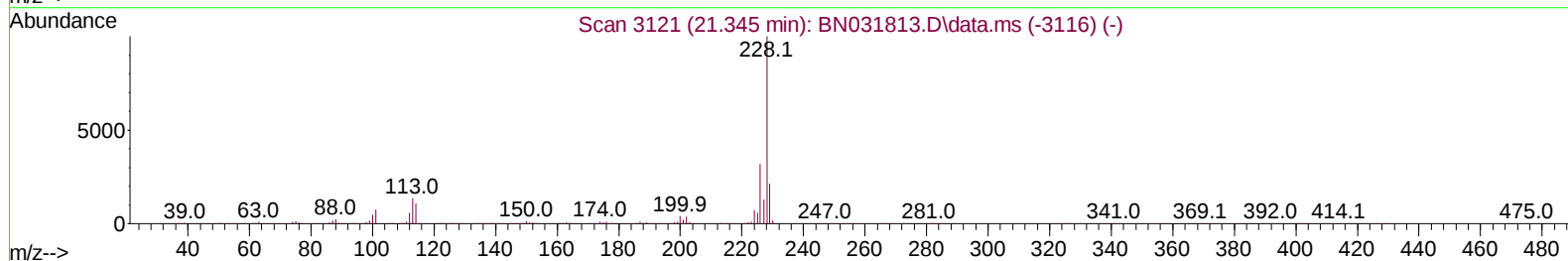
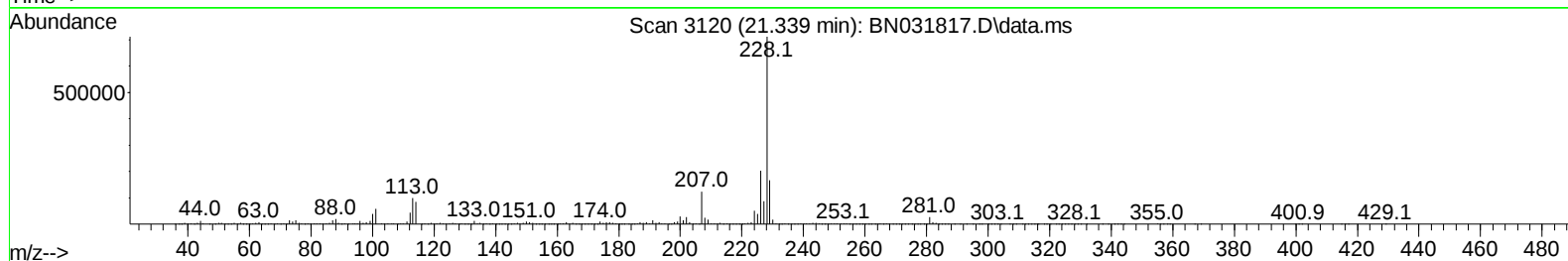
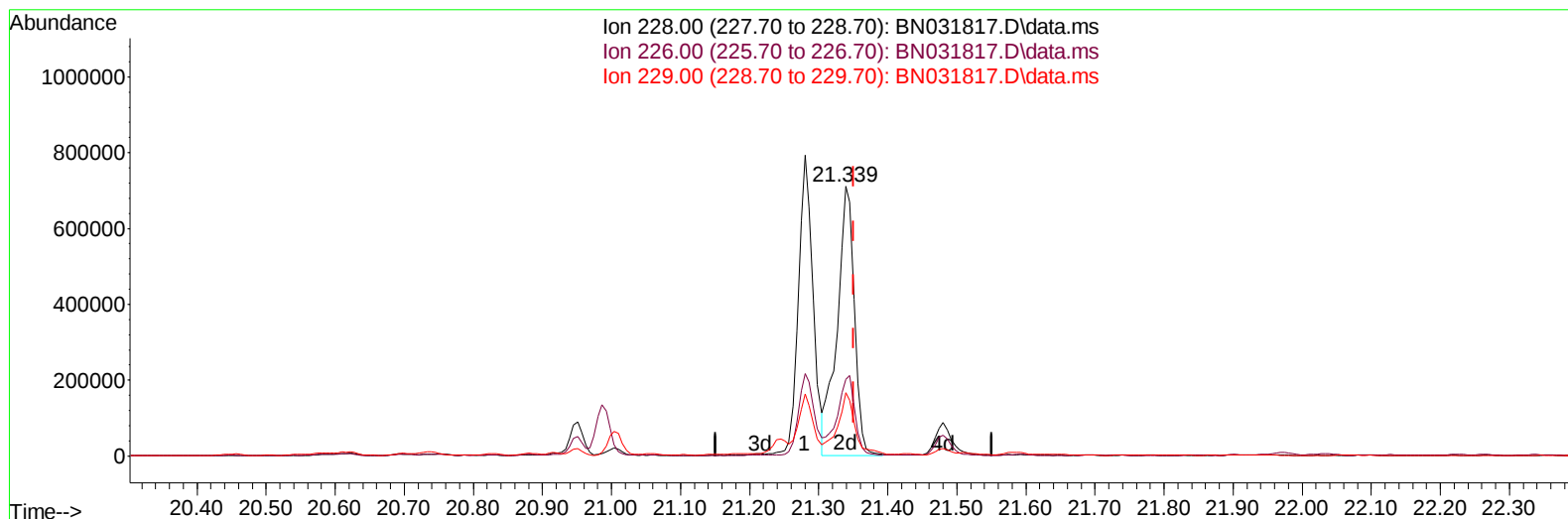
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TIC: BN031817.D\data.ms

(87) Chrysene

21.339min (-0.012) 13.82 ng/ul m

response 1247551

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	29.60	28.54
229.00	18.90	23.40#
0.00	0.00	0.00

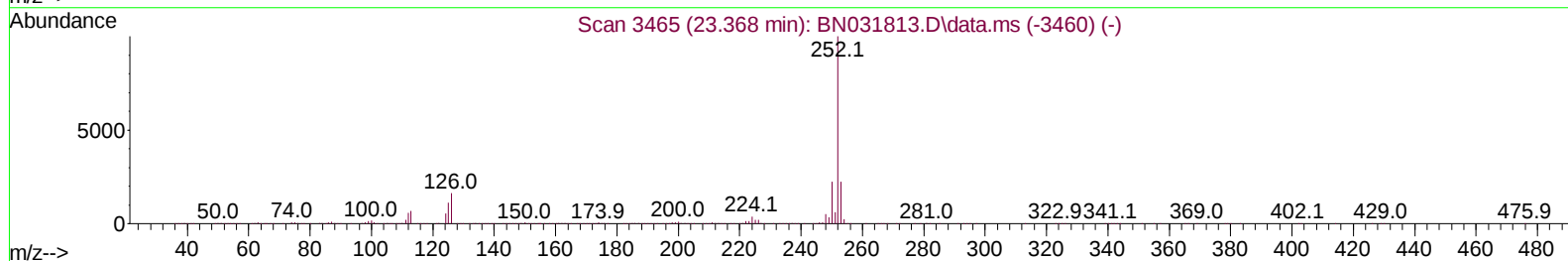
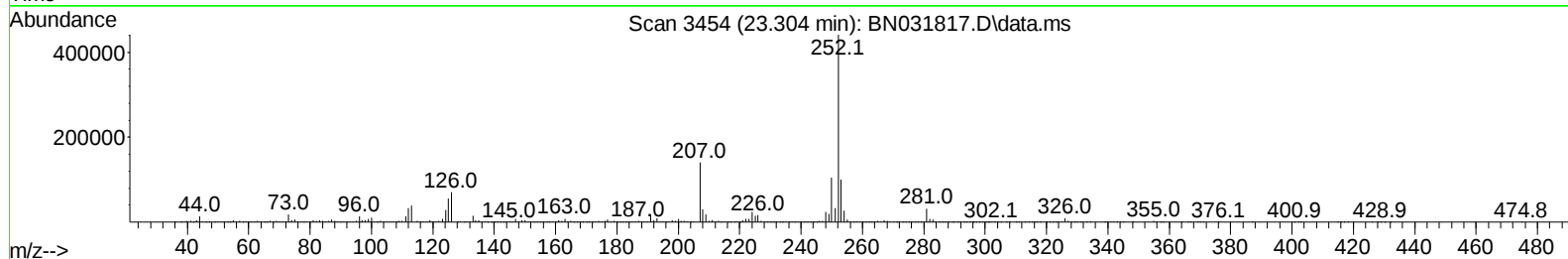
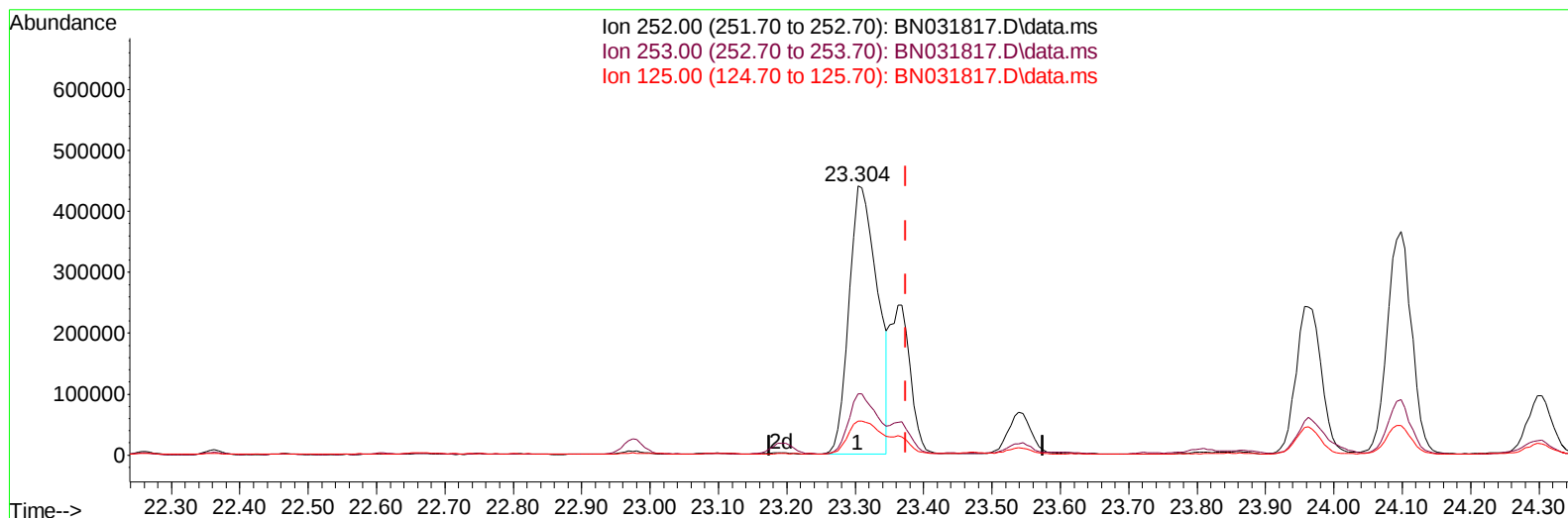
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Operator : MA/JU
Sample : P2634-20
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHHY4

Manual IntegrationsAPPROVED

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TIC: BN031817.D\data.ms

(91) Benzo(k)fluoranthene

23.304min (-0.071) 17.71 ng/u1

response 1273538

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.00	22.85
125.00	11.80	12.67
0.00	0.00	0.00

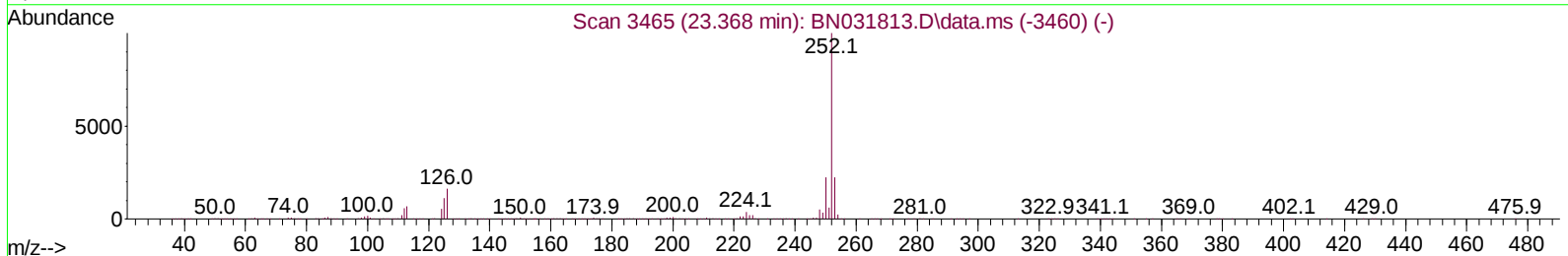
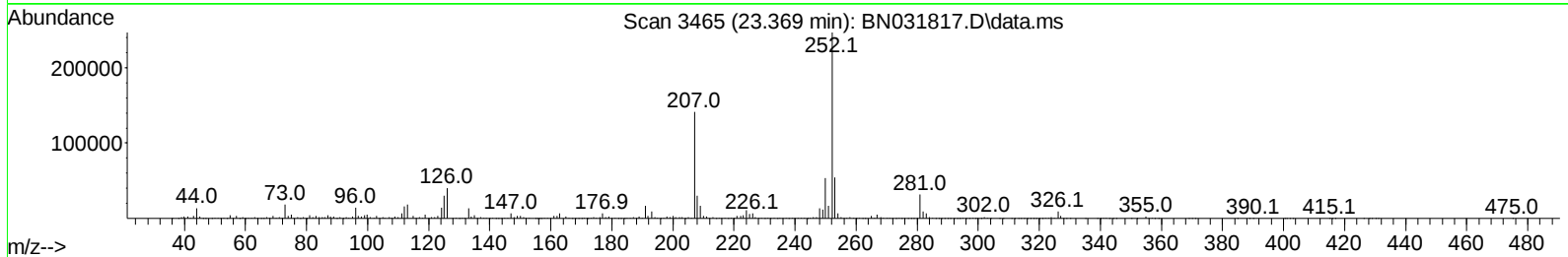
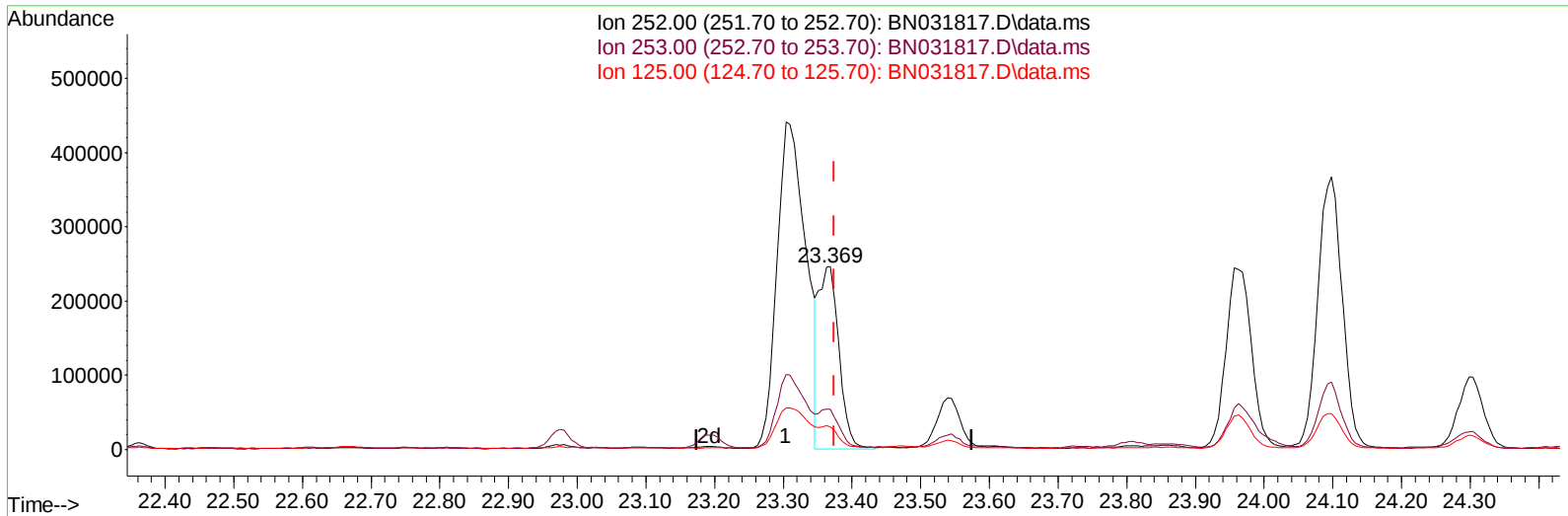
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ALS Vial : 6 Sample Multiplier: 1

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TIC: BN031817.D\data.ms

(91) Benzo(k)fluoranthene

23.369min (-0.006) 7.02 ng/u1 m

response 505054

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.00	22.04
125.00	11.80	12.16
0.00	0.00	0.00

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 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHBY4

Manual IntegrationsAPPROVED

Quant Time: Jun 06 14:36:45 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 29 15:13:07 2024
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.669	152		314480	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.457	136		1321233	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.316	164		806292	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.063	188		1644039	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240		1444021	20.000	ng/ul	0.00
88) Perylene-d12	24.233	264		1226279	20.000	ng/ul	-0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.187	96		44196	5.814	ng/uL	0.00
4) Pyridine-d5	3.587	84		552403	25.815	ng/ul	-0.01
7) Phenol-d5	6.846	99		798139	29.605	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.011	67		460008	29.171	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.205	132		647534	31.478	ng/ul	-0.01
15) 4-Methylphenol-d8	8.375	113		607760	27.655	ng/ul	-0.01
21) Nitrobenzene-d5	8.828	128		324644	32.136	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.546	143		339606	32.544	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.075	165		587519	30.157	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.593	131		790398	27.325	ng/ul	-0.02
46) Dimethylphthalate-d6	13.728	166		1701194	30.286	ng/ul	-0.02
49) Acenaphthylene-d8	14.004	160		2059504	31.887	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.522	143		306997	27.297	ng/ul	0.00
60) Fluorene-d10	15.310	176		1506025	31.687	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.440	200		217454	26.314	ng/ul	0.00
73) Anthracene-d10	17.163	188		2270126	32.253	ng/ul	0.00
81) Pyrene-d10	19.463	212		2662928	34.461	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.033	264		1881840	32.138	ng/ul	0.00
Target Compounds							
						Qvalue	
50) Acenaphthylene	14.034	152		83289	1.141	ng/ul	98
52) Acenaphthene	14.381	153		73052	1.513	ng/ul	97
61) Fluorene	15.369	166		59073	1.118	ng/ul	96
72) Phenanthrene	17.104	178		1343818	16.046	ng/ul	99
74) Anthracene	17.198	178		234341	2.761	ng/ul	100
77) Carbazole	17.469	167		131386	1.816	ng/ul	96
80) Fluoranthene	19.128	202		2961821	32.072	ng/ul	100
82) Pyrene	19.492	202		2716412	28.669	ng/ul	98
85) Benzo(a)anthracene	21.280	228		1163391	12.037	ng/ul	99
87) Chrysene	21.339	228		1247551m	13.818	ng/ul	
90) Benzo(b)fluoranthene	23.304	252		1273538	17.375	ng/ul	97
91) Benzo(k)fluoranthene	23.369	252		505054m	7.022	ng/ul	
93) Benzo(a)pyrene	24.098	252		908771	13.238	ng/ul	94
94) Indeno(1,2,3-cd)pyrene	27.533	276		635534	8.034	ng/ul	98
95) Dibenzo(a,h)anthracene	27.568	278		143635	2.224	ng/ul#	91
96) Benzo(g,h,i)perylene	28.568	276		620518	9.868	ng/ul	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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