

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122022\
 Data File : BM038174.D
 Acq On : 21 Dec 2022 16:25
 Operator : CG/JU
 Sample : N6014-21
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBXT3

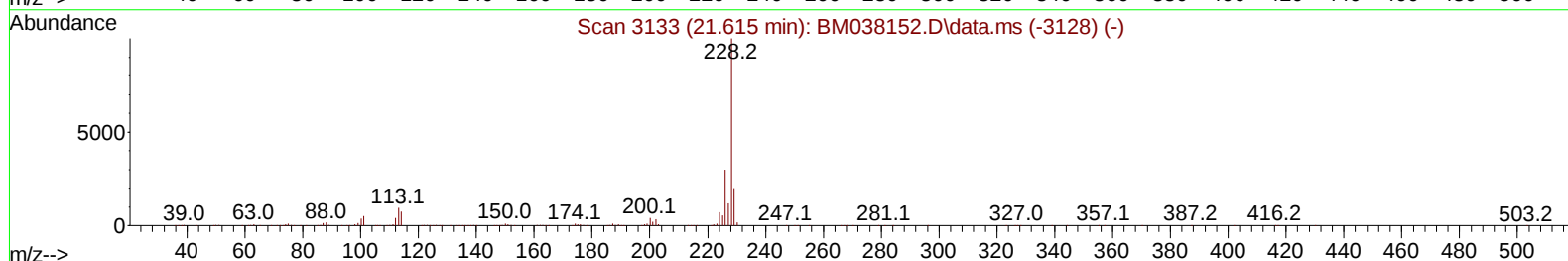
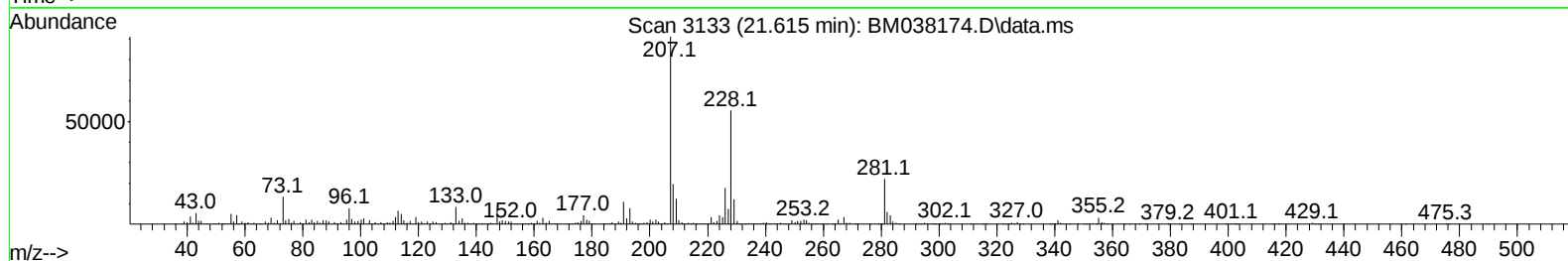
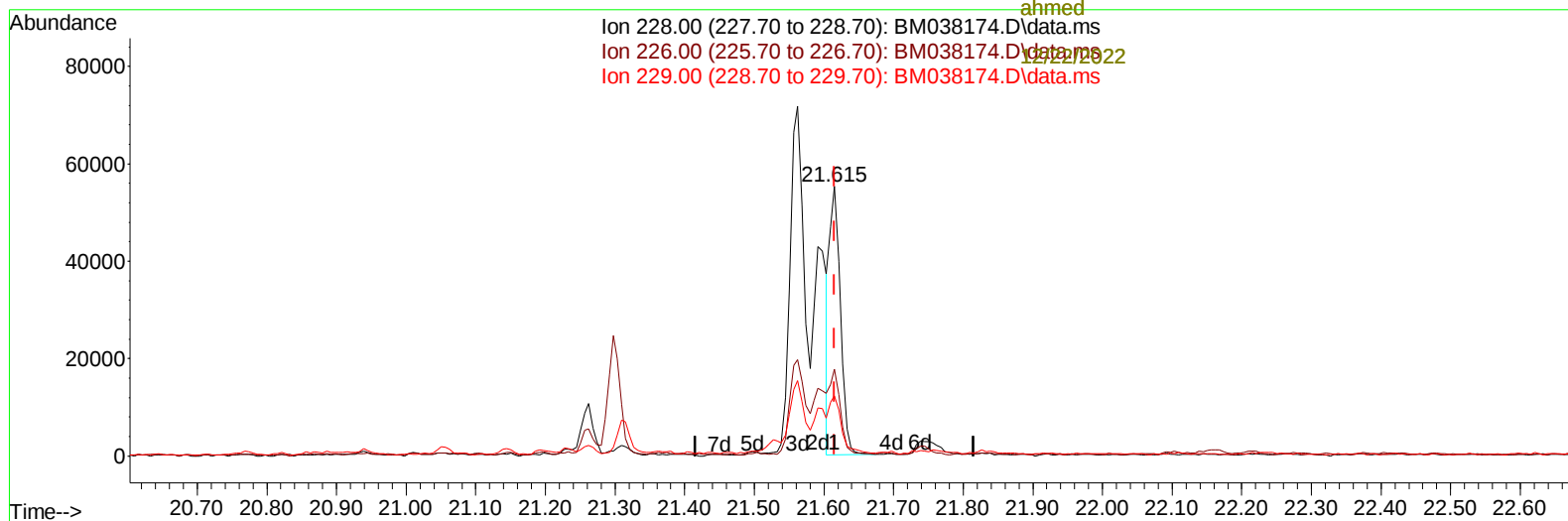
Manual IntegrationsAPPROVED

Quant Time: Dec 21 23:32:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 21 00:11:56 2022
 Response via : Initial Calibration

Reviewed By :Jagrut
 Upadhyay

12/22/2022
 Supervised By :mohammad
 ahmed

Ion 228.00 (227.70 to 228.70): BM038174.D\data.ms
 Ion 226.00 (225.70 to 226.70): BM038174.D\data.ms
 Ion 229.00 (228.70 to 229.70): BM038174.D\data.ms



TIC: BM038174.D\data.ms

(87) Chrysene

21.615min (-0.000) 1.27 ng/u1

response 59500

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	30.40	32.18
229.00	19.50	22.29
0.00	0.00	0.00

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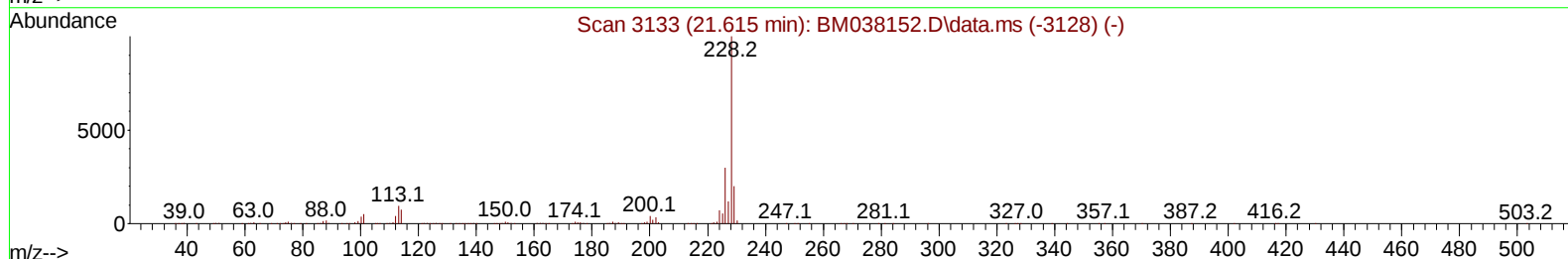
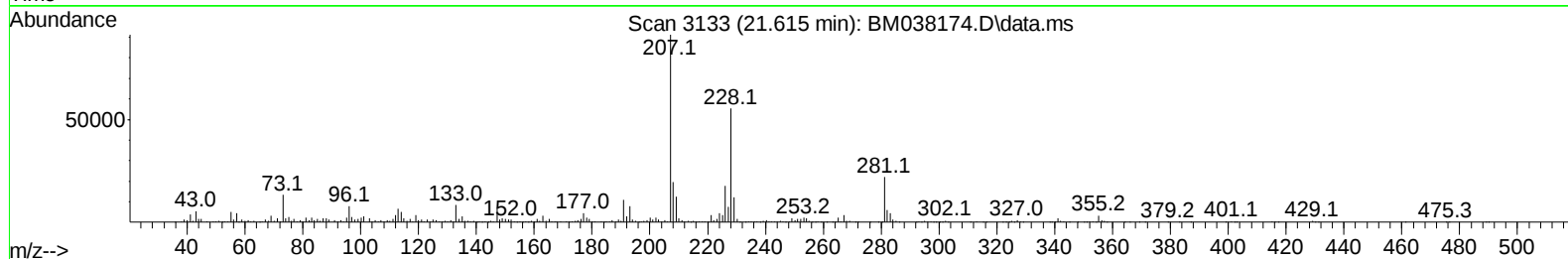
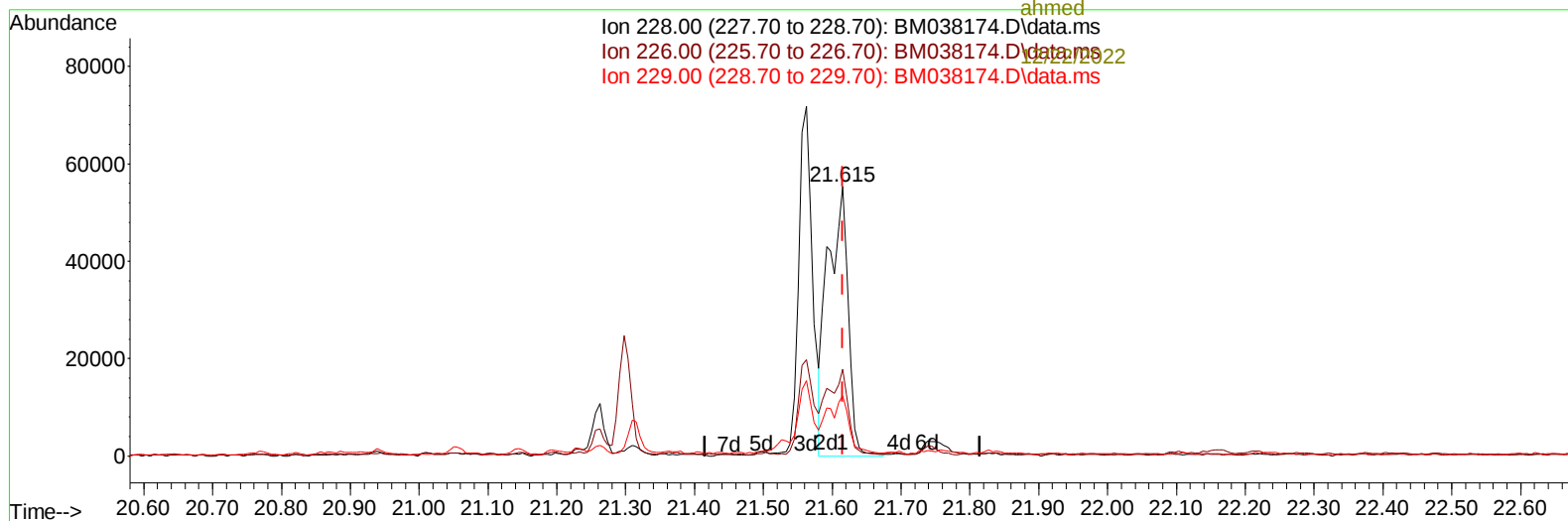
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TIC: BM038174.D\data.ms

(87) Chrysene

21.615min (-0.000) 2.45 ng/ul m

response 114627

Ion	Exp%	Act%
228.00	100.00	100.00
226.00	30.40	32.18
229.00	19.50	22.29
0.00	0.00	0.00

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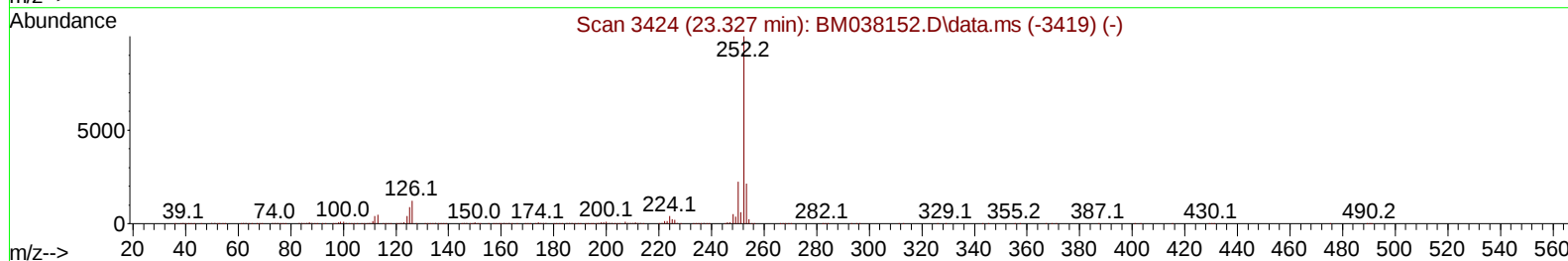
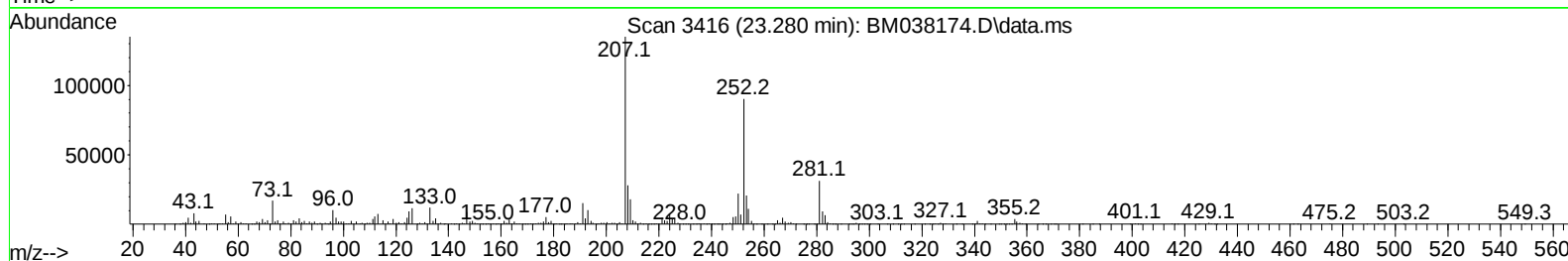
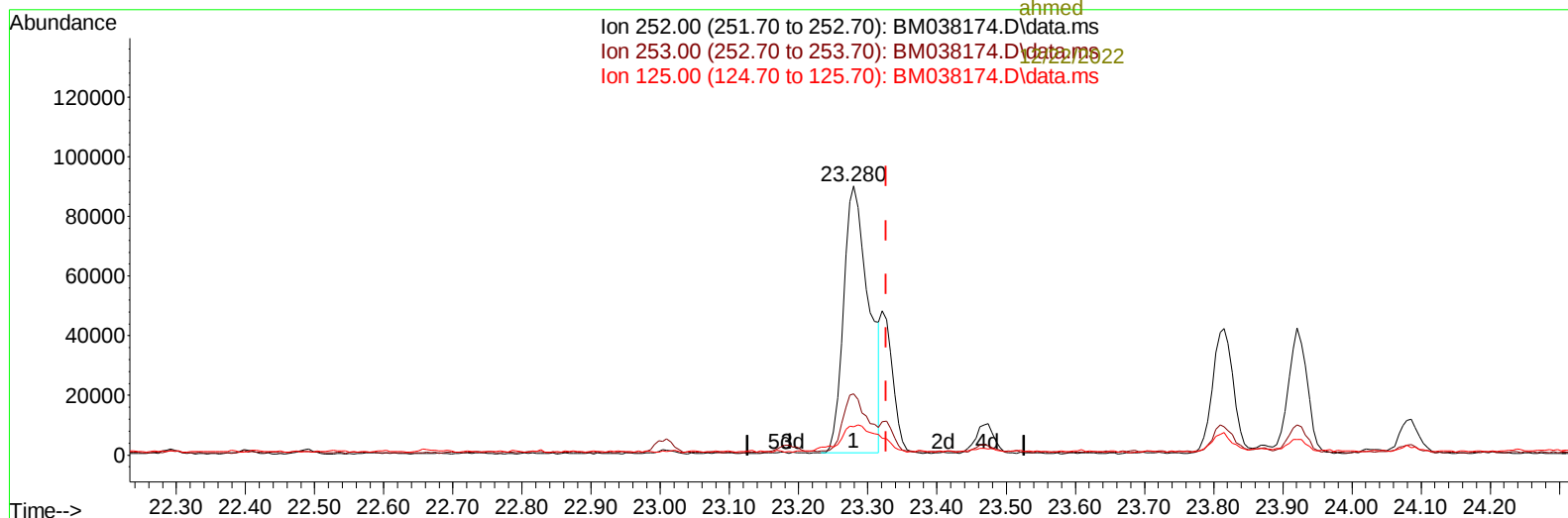
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Ion 252.00 (251.70 to 252.70): BM038174.D\data.ms
 Ion 253.00 (252.70 to 253.70): BM038174.D\data.ms
 Ion 125.00 (124.70 to 125.70): BM038174.D\data.ms



TIC: BM038174.D\data.ms

(91) Benzo(k)fluoranthene

23.280min (-0.047) 4.87 ng/u1

response 227323

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.80	22.81
125.00	10.10	10.45
0.00	0.00	0.00

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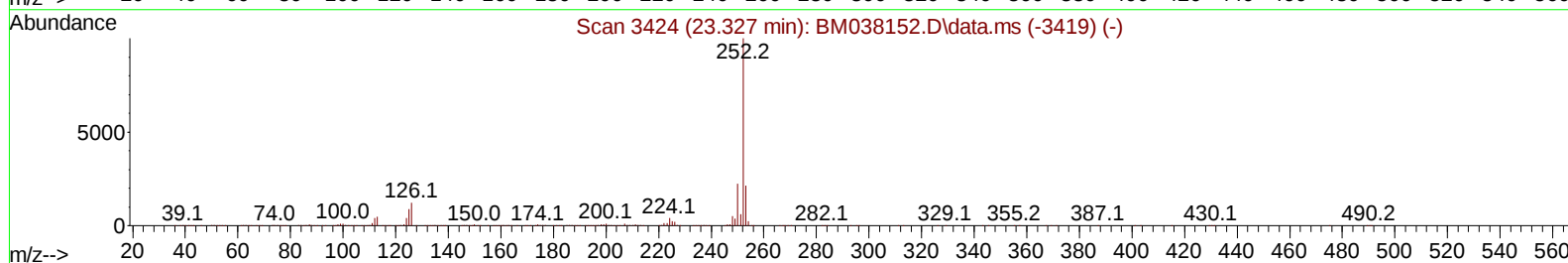
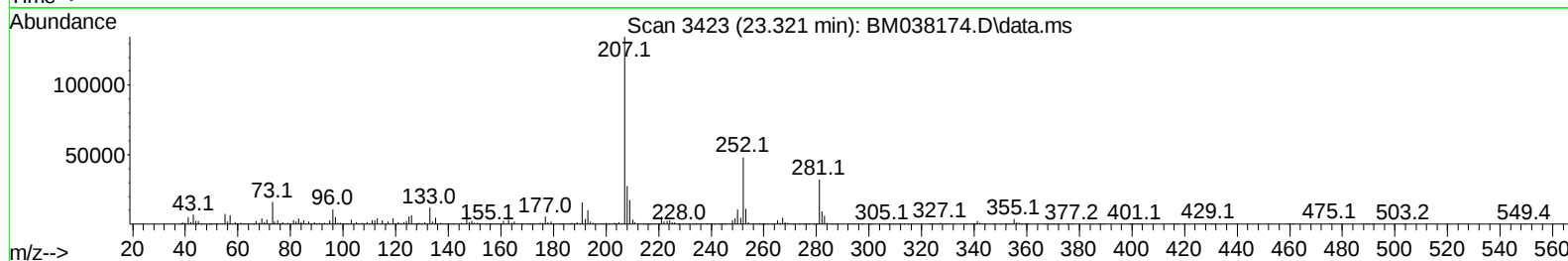
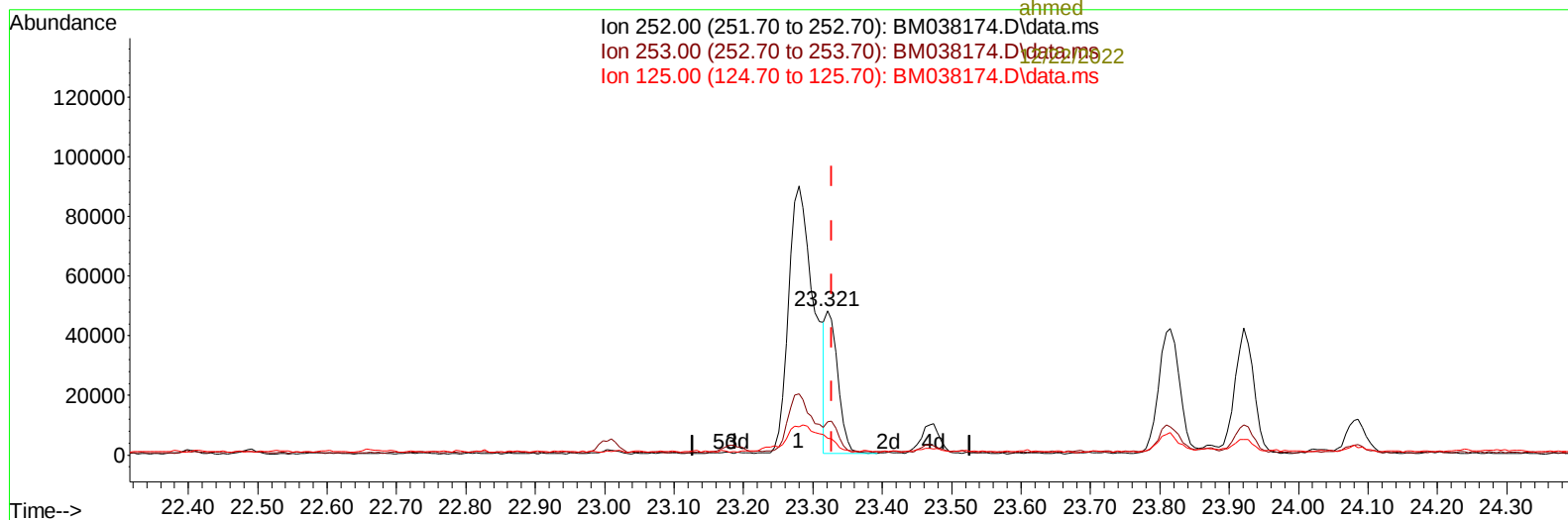
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TIC: BM038174.D\data.ms

(91) Benzo(k)fluoranthene

23.321min (-0.006) 1.24 ng/ul m

response 57918

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.80	23.17
125.00	10.10	11.50
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----12/22/2022						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.040	152	214530	20.000	ng/ul	0.00
20) Naphthalene-d8	10.857	136	900238	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.669	164	534277	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.410	188	1064120	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.580	240	739766	20.000	ng/ul	0.00
88) Perylene-d12	24.033	264	776361	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	19008	4.023	ng/uL	0.00
4) Pyridine-d5	3.828	84	312427	22.291	ng/ul	0.00
7) Phenol-d5	7.204	99	581718	33.369	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.369	67	279796	26.358	ng/ul	0.00
11) 2-Chlorophenol-d4	7.569	132	495419	34.643	ng/ul	0.00
15) 4-Methylphenol-d8	8.757	113	494492	36.339	ng/ul	0.00
21) Nitrobenzene-d5	9.216	128	250093	35.087	ng/ul	0.00
24) 2-Nitrophenol-d4	9.939	143	299960	40.158	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.475	165	582942	41.088	ng/ul	0.00
31) 4-Chloroaniline-d4	10.998	131	606749	30.982	ng/ul	0.00
46) Dimethylphthalate-d6	14.080	166	1653719	43.416	ng/ul	0.00
49) Acenaphthylene-d8	14.369	160	1920683	41.258	ng/ul	0.00
54) 4-Nitrophenol-d4	14.874	143	245636	37.839	ng/ul	0.00
60) Fluorene-d10	15.657	176	1446199	42.616	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.780	200	129940	22.840	ng/ul	0.00
73) Anthracene-d10	17.510	188	2090116	42.085	ng/ul	0.00
81) Pyrene-d10	19.792	212	2075724	46.711	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.874	264	1631581	42.075	ng/ul	0.00
Target Compounds						
					Qvalue	
80) Fluoranthene	19.456	202	68886	1.313	ng/ul	98
82) Pyrene	19.821	202	211469	3.874	ng/ul	96
85) Benzo(a)anthracene	21.562	228	99492	1.998	ng/ul	98
87) Chrysene	21.615	228	114627m	2.453	ng/ul	
90) Benzo(b)fluoranthene	23.280	252	227323	4.723	ng/ul	98
91) Benzo(k)fluoranthene	23.321	252	57918m	1.242	ng/ul	
93) Benzo(a)pyrene	23.921	252	79816	1.886	ng/ul	96

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

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