

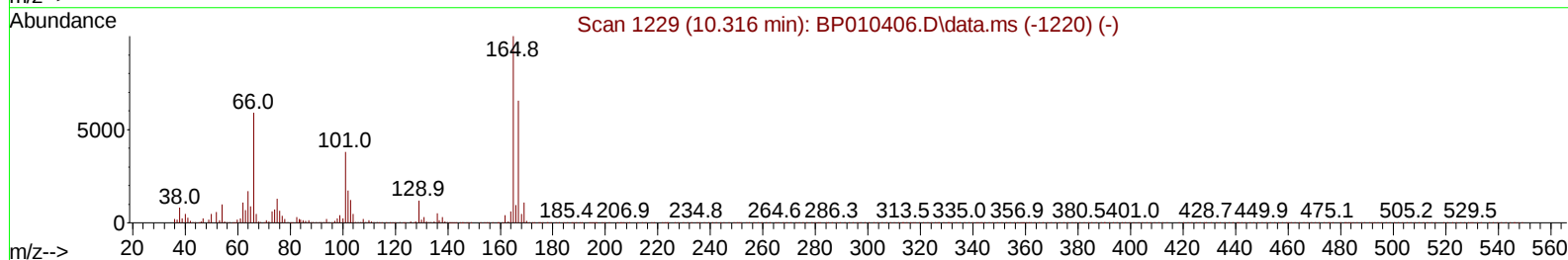
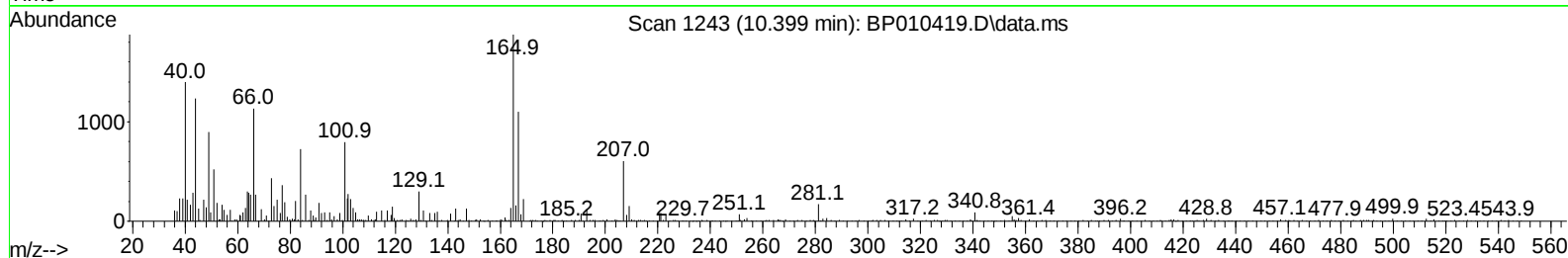
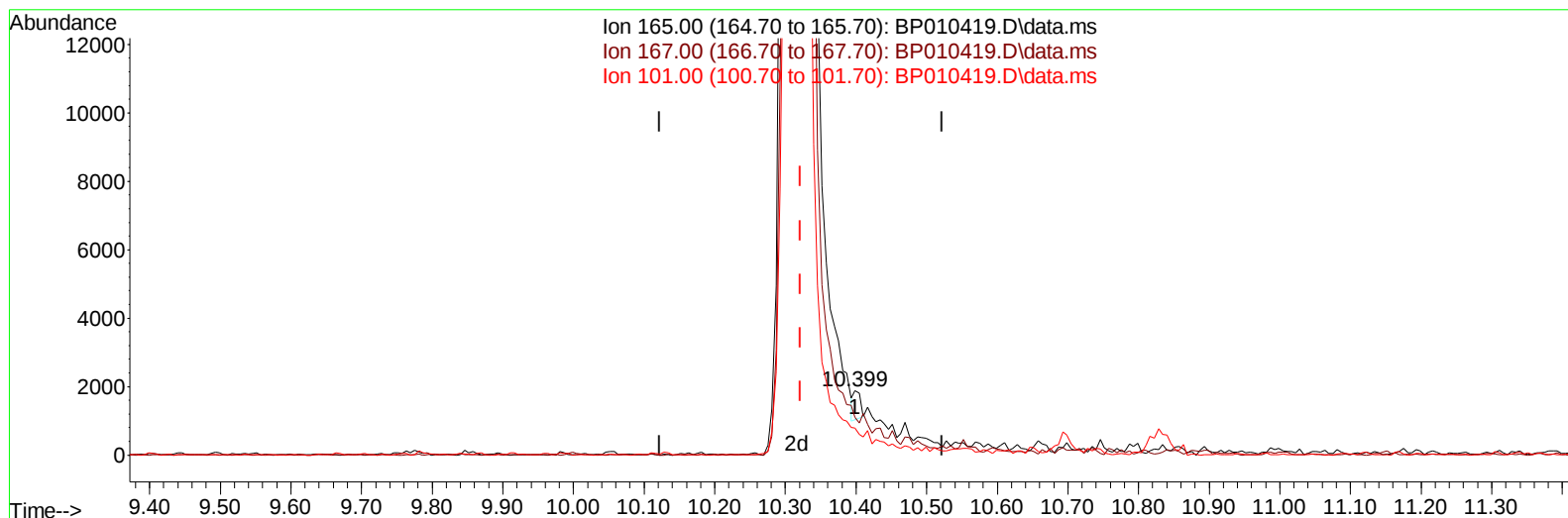
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051922\
Data File : BP010419.D
Acq On : 20 May 2022 00:35
Operator : CG/JU
Sample : N2918-08
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTF3

Manual IntegrationsAPPROVED

Quant Time: May 20 02:29:10 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 13 15:04:31 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/20/2022
Supervised By :Yogesh Patel 05/24/2022



TIC: BP010419.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.399min (+ 0.077) 0.05 ng/u1

response 581

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	58.64
101.00	49.80	42.48
0.00	0.00	0.00

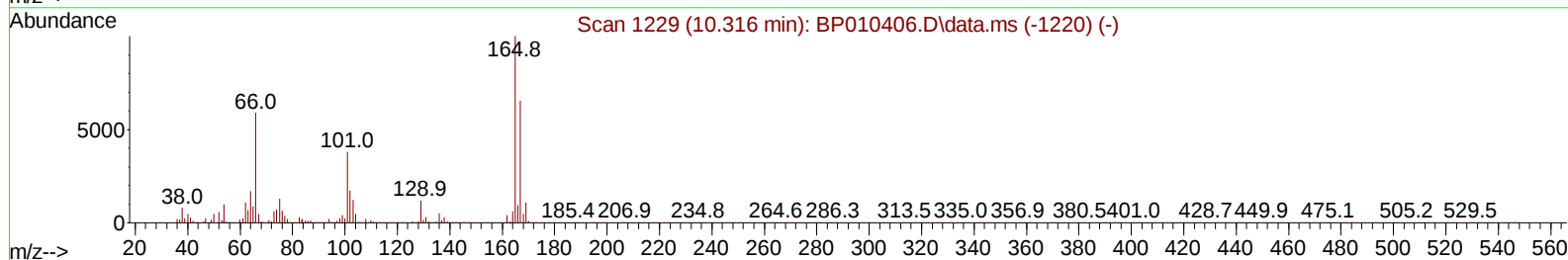
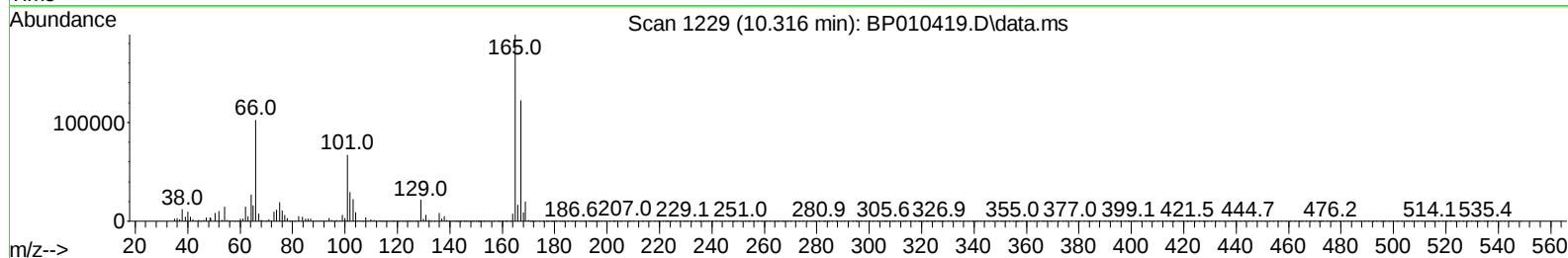
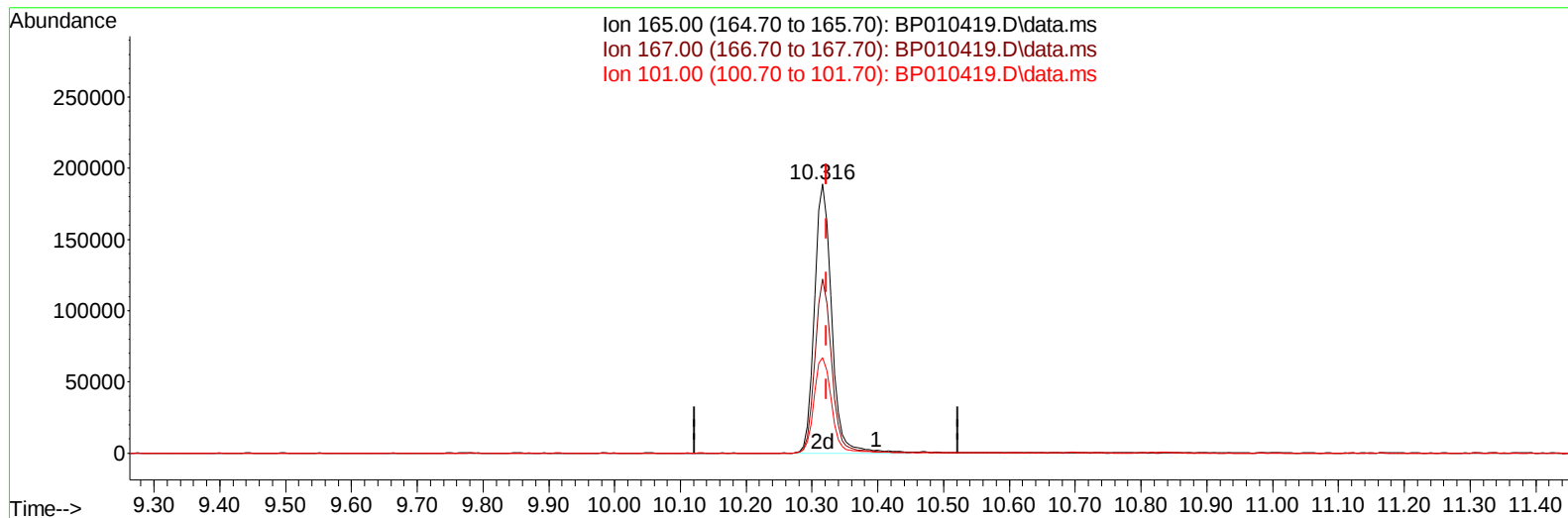
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TIC: BP010419.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.316min (-0.006) 29.47 ng/ul m

response 339932

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	64.67
101.00	49.80	35.62#
0.00	0.00	0.00

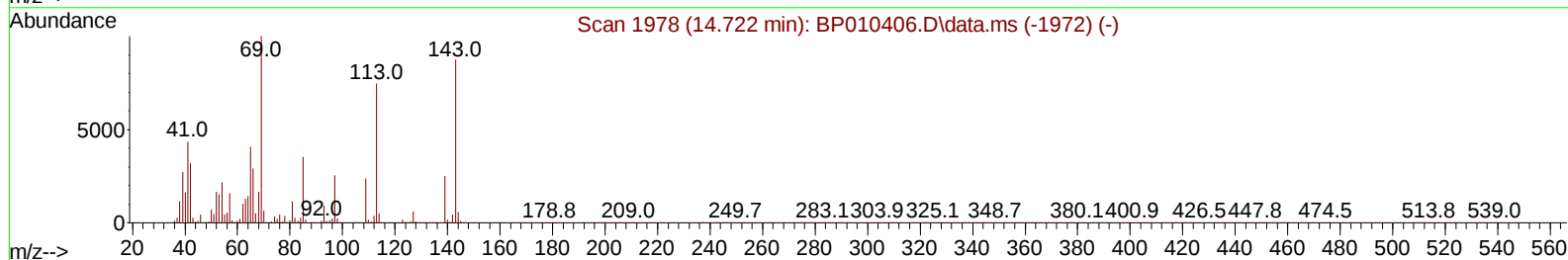
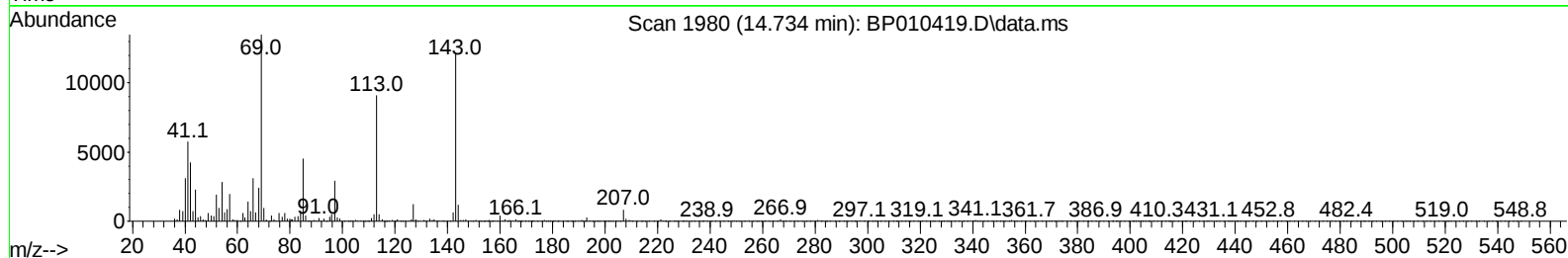
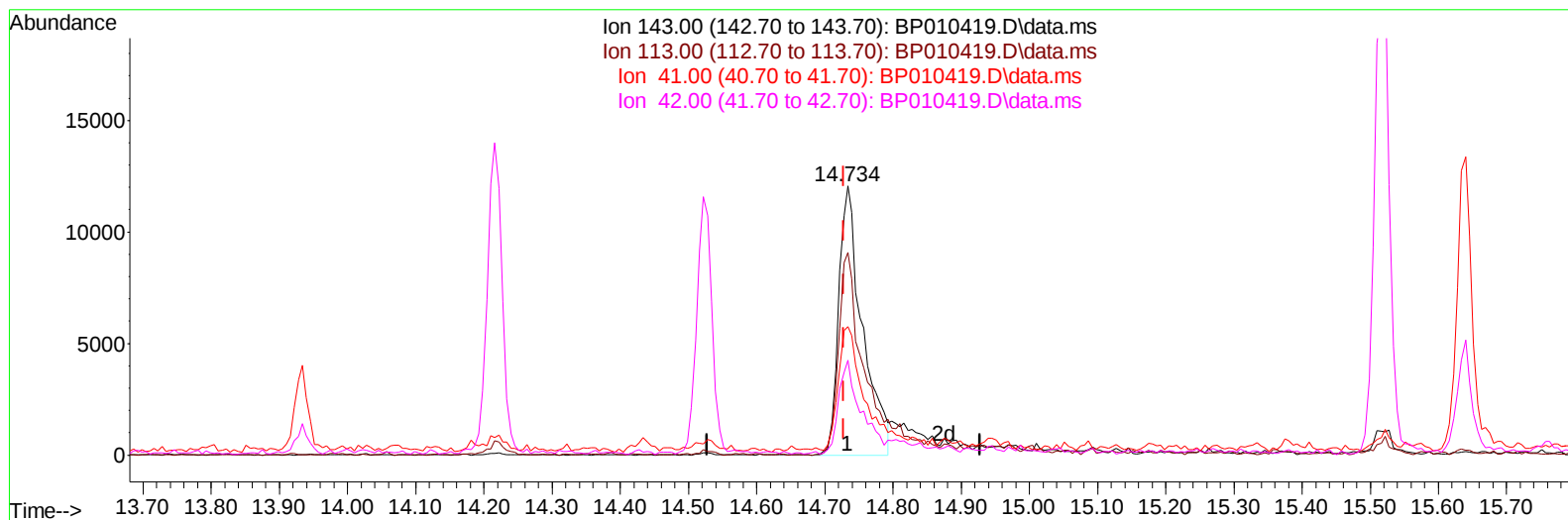
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
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TIC: BP010419.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 4.34 ng/ul

response 29177

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	75.45#
41.00	23.20	47.78#
42.00	18.00	35.19#

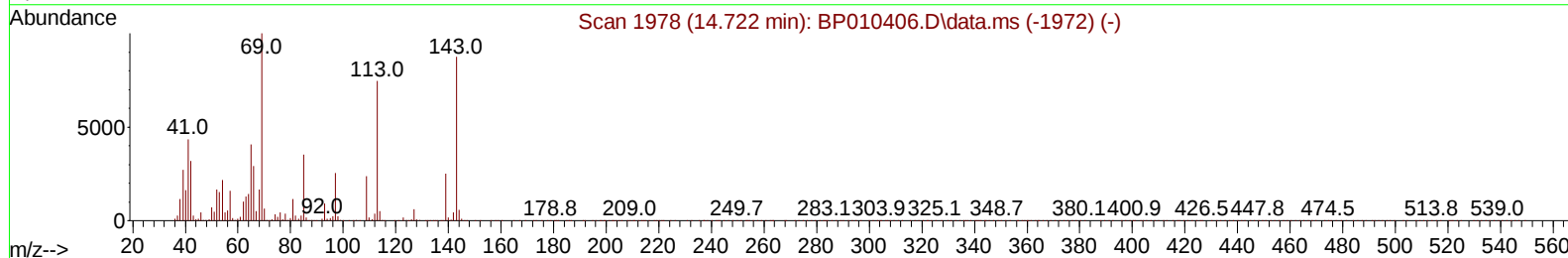
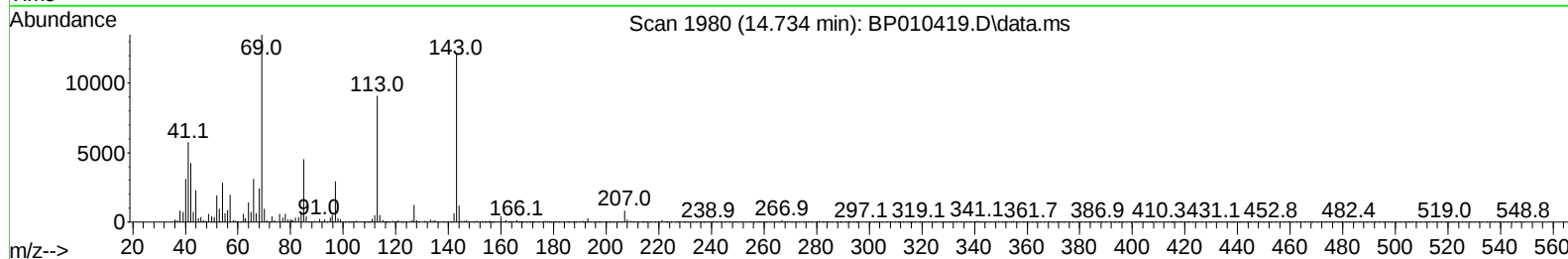
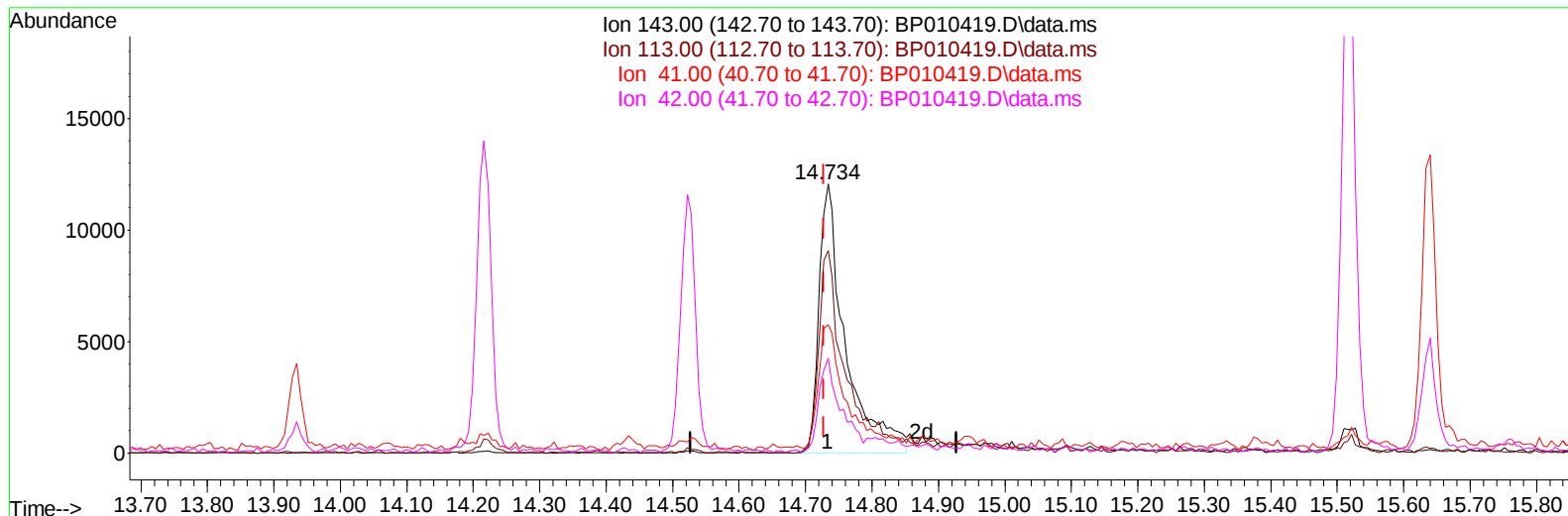
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP010419.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 4.94 ng/ul m

response 33209

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	75.45#
41.00	23.20	47.78#
42.00	18.00	35.19#

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 Operator : CG/JU
 Sample : N2918-08
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBTF3

Manual IntegrationsAPPROVED

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 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	167590	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	743684	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.522	164	512776	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.275	188	1094123	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.363	240	989955	20.000	ng/ul	# 0.00
88) Perylene-d12	23.786	264	844805	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	17272	4.393	ng/uL	0.00
4) Pyridine-d5	3.770	84	97971	8.776	ng/ul	0.00
7) Phenol-d5	7.058	99	89746	6.380	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	315585	33.393	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132	276516	25.690	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	190699	15.953	ng/ul	-0.01
21) Nitrobenzene-d5	9.052	128	198766	34.534	ng/ul	0.00
24) 2-Nitrophenol-d4	9.775	143	207603	34.036	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	339932m	29.474	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	413485	24.112	ng/ul	-0.01
46) Dimethylphthalate-d6	13.934	166	1344232	35.581	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1514664	34.758	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	33209m	4.939	ng/ul	0.00
60) Fluorene-d10	15.516	176	1183836	36.018	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	207624	32.447	ng/ul	0.00
73) Anthracene-d10	17.375	188	2024998	40.873	ng/ul	0.00
81) Pyrene-d10	19.610	212	2383911	45.344	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.633	264	1812349	42.723	ng/ul	0.00

Target Compounds Qvalue

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

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