

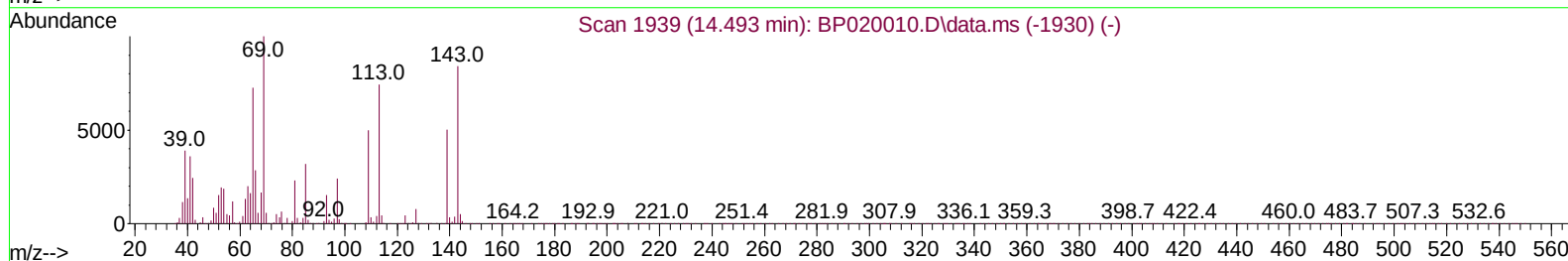
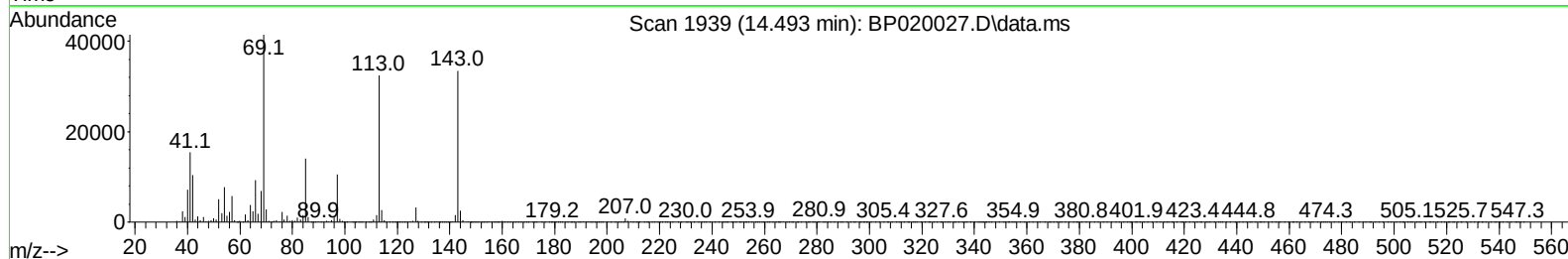
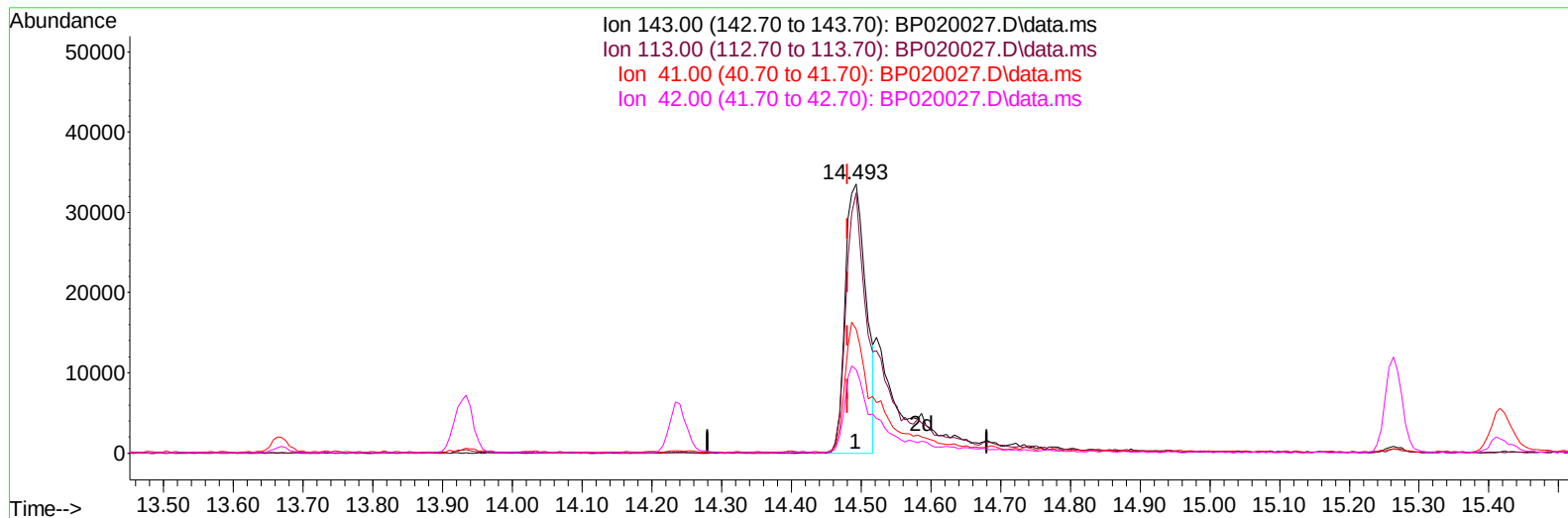
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032824\
 Data File : BP020027.D
 Acq On : 29 Mar 2024 02:40
 Operator : MA/JU
 Sample : PB159810TB 10X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLEB810

Manual IntegrationsAPPROVED

Quant Time: Mar 29 04:43:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 28 14:08:48 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 03/29/2024
 Supervised By :Jagrut Upadhyay 03/29/2024



TIC: BP020027.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.493min (+ 0.012) 18.64 ng/ul

response 69388

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	96.68
41.00	40.90	46.09
42.00	28.50	31.09

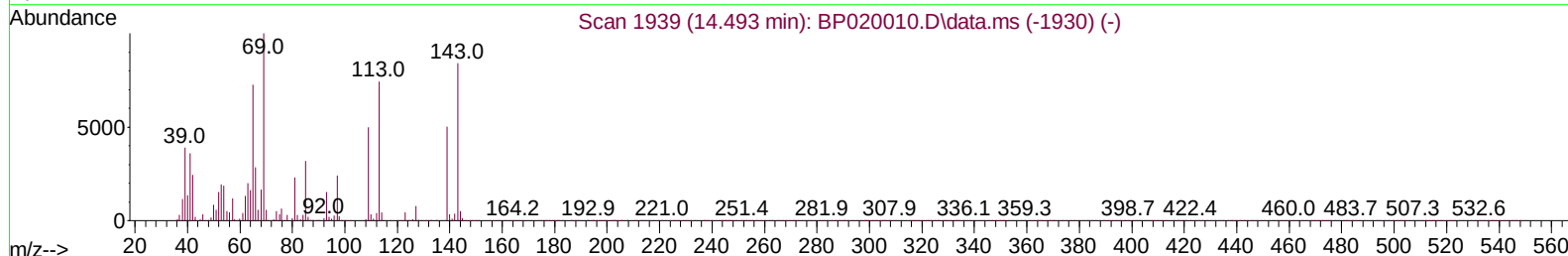
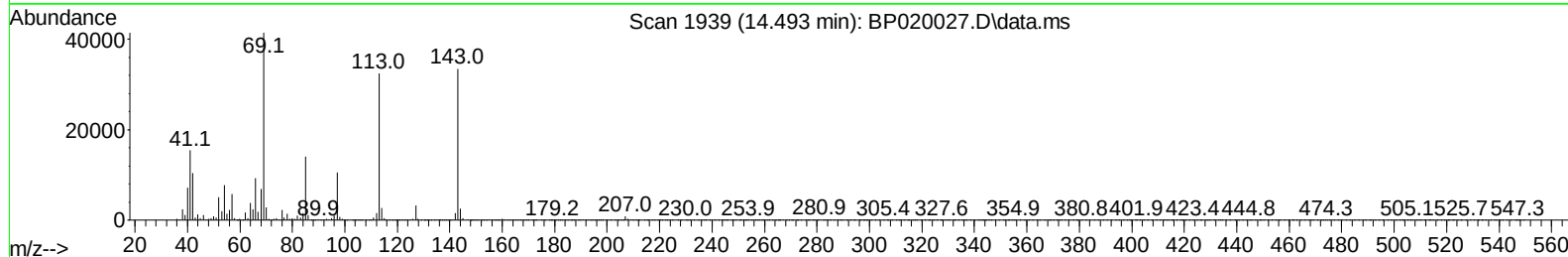
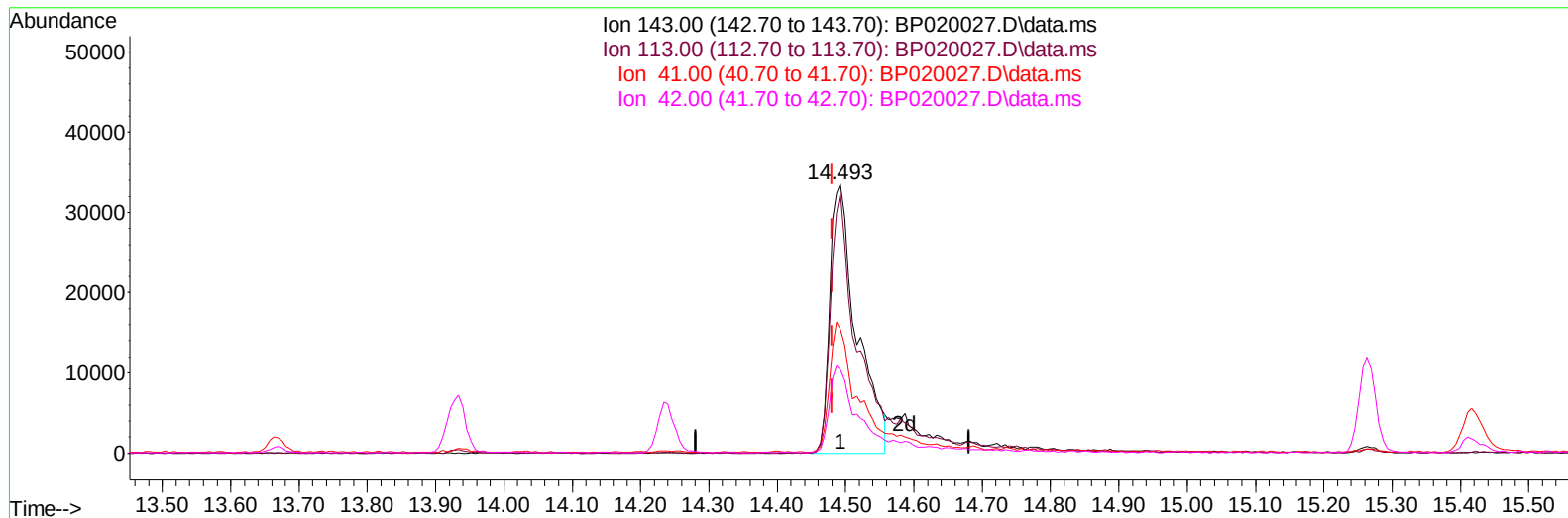
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(54) 4-Nitrophenol-d4 (S)

14.493min (+ 0.012) 24.63 ng/ul m

response 91678

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	96.68
41.00	40.90	46.09
42.00	28.50	31.09

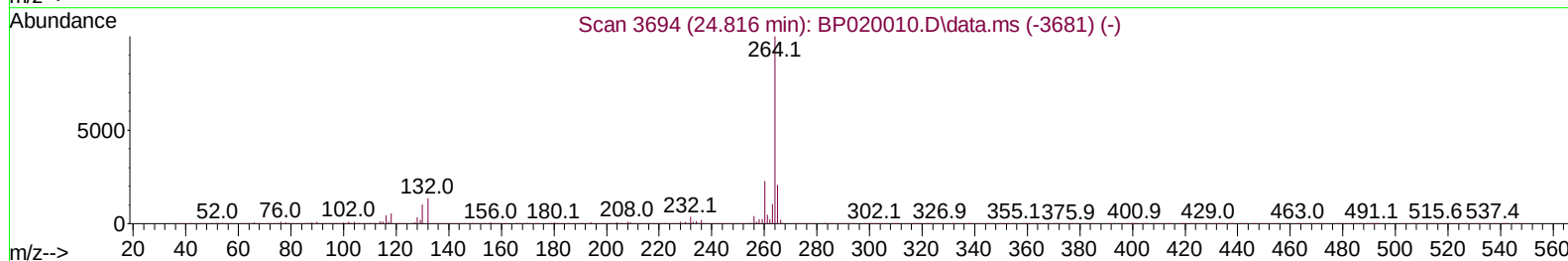
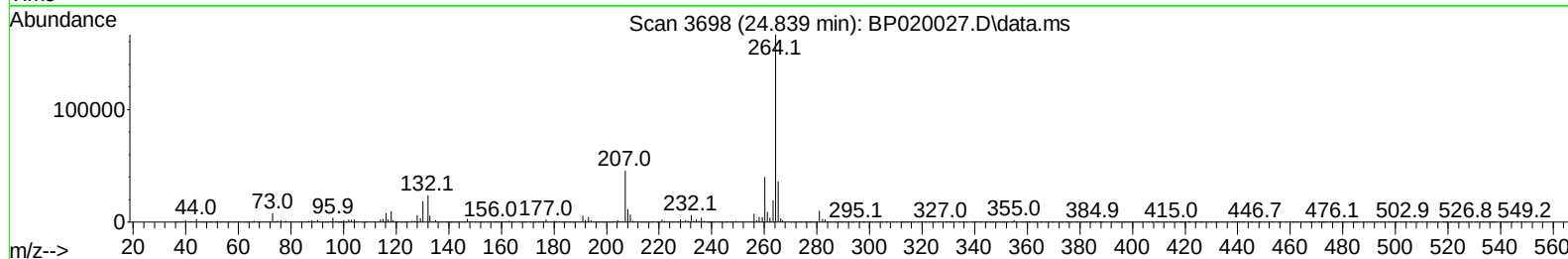
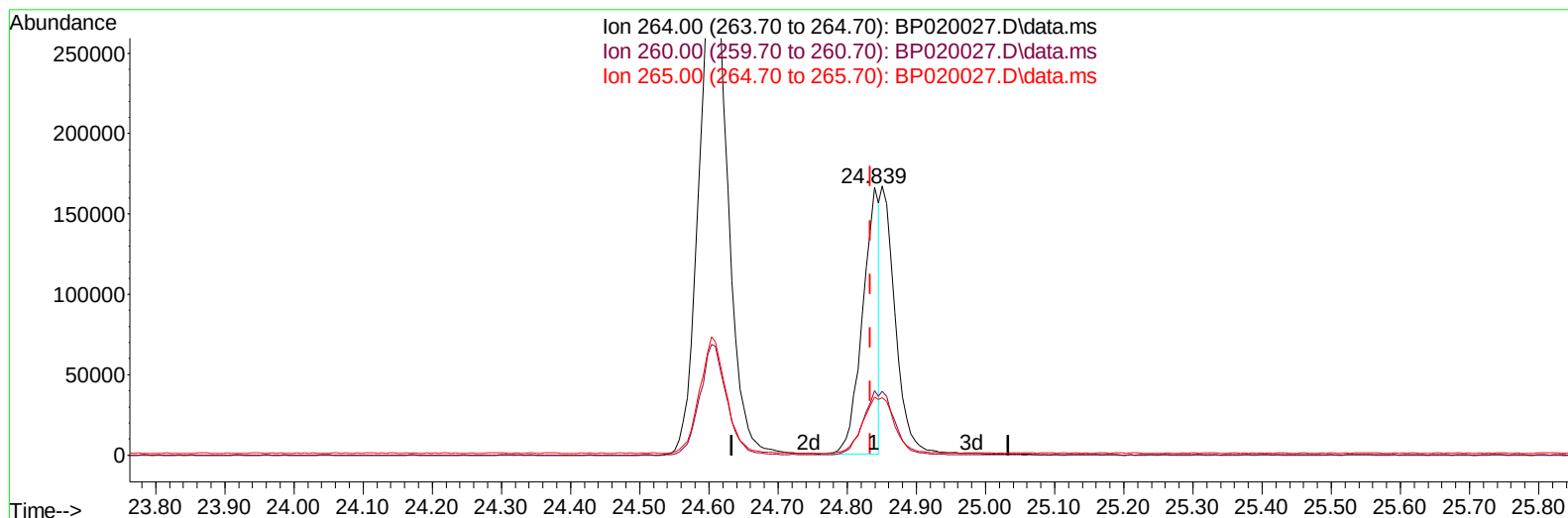
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TIC: BP020027.D\data.ms

(88) Perylene-d12 (I)

24.839min (+ 0.006) 20.00 ng/ul

response 277467

Ion	Exp%	Act%
264.00	100.00	100.00
260.00	24.30	24.12
265.00	21.80	21.65
0.00	0.00	0.00

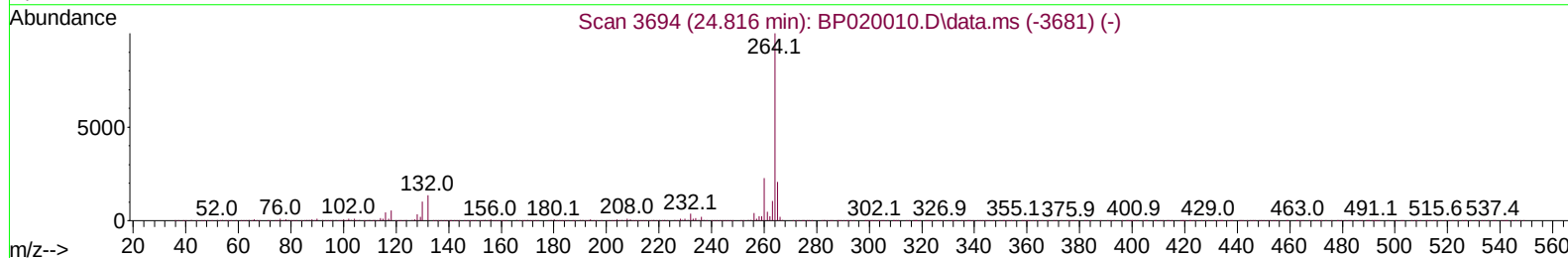
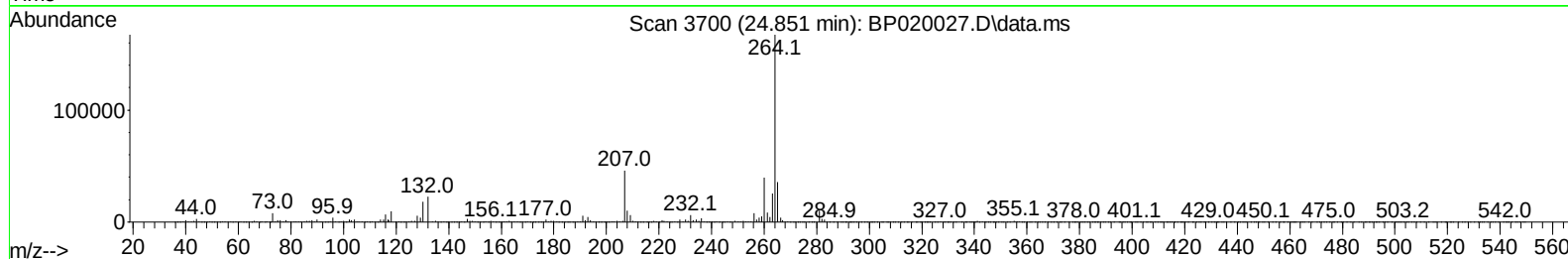
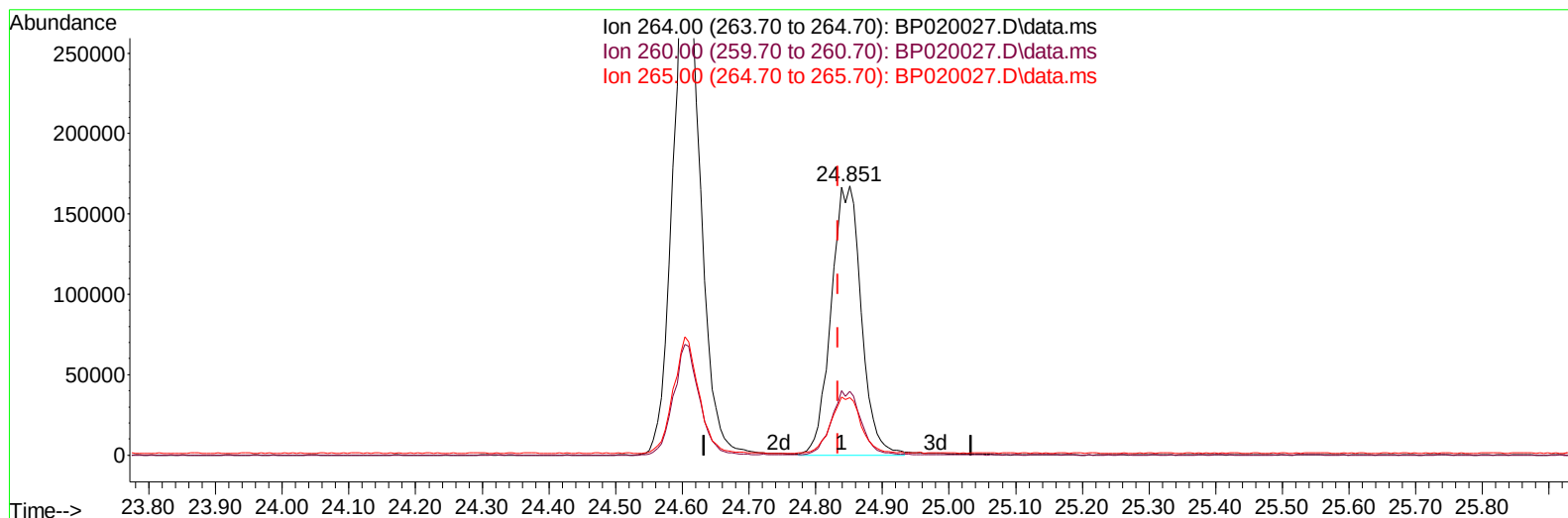
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Instrument :
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TIC: BP020027.D\data.ms

(88) Perylene-d12 (I)

24.851min (+ 0.018) 20.00 ng/ul m

response 526168

Ion	Exp%	Act%
264.00	100.00	100.00
260.00	24.30	23.71
265.00	21.80	21.31
0.00	0.00	0.00

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 ALS Vial : 18 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 03/29/2024
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Quant Time: Mar 29 04:45:56 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.611	152	116142	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	494797	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.240	164	337711	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.087	188	673851	20.000	ng/ul	0.01
79) Chrysene-d12	21.545	240	503534	20.000	ng/ul	0.00
88) Perylene-d12	24.851	264	526168m	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	21451	8.092	ng/uL	0.00
4) Pyridine-d5	3.640	84	182069	22.747	ng/ul	0.00
7) Phenol-d5	6.793	99	349505	35.681	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.964	67	246872	38.684	ng/ul	0.00
11) 2-Chlorophenol-d4	7.152	132	261632	37.192	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	260489	34.373	ng/ul	0.00
21) Nitrobenzene-d5	8.758	128	137076	36.144	ng/ul	0.00
24) 2-Nitrophenol-d4	9.463	143	153134	36.089	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.999	165	262217	33.837	ng/ul	0.00
31) 4-Chloroaniline-d4	10.552	131	19789	1.959	ng/ul	0.04
46) Dimethylphthalate-d6	13.669	166	912891	38.555	ng/ul	0.02
49) Acenaphthylene-d8	13.934	160	973790	35.207	ng/ul	0.00
54) 4-Nitrophenol-d4	14.493	143	91678m	24.629	ng/ul	0.01
60) Fluorene-d10	15.263	176	754829	37.423	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.416	200	132043	32.736	ng/ul	0.00
73) Anthracene-d10	17.175	188	1118009	37.481	ng/ul	0.00
81) Pyrene-d10	19.575	212	1198132	40.555	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.604	264	927010	36.672	ng/ul	0.00

Target Compounds Qvalue

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

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