

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071423\
 Data File : BP016237.D
 Acq On : 14 Jul 2023 21:22
 Operator : MA/JU
 Sample : 03457-09
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0C51

Quant Time: Jul 14 23:57:22 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 14 23:53:26 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	126388	20.000	ng/ul	0.00
20) Naphthalene-d8	10.786	136	614638	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.616	164	403805	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.380	188	926281	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	1003734	20.000	ng/ul	0.01
88) Perylene-d12	23.968	264	1165617	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	8507	2.382	ng/uL	0.00
4) Pyridine-d5	3.746	84	316	0.035	ng/ul	0.00
7) Phenol-d5	7.181	99	120069	10.518	ng/ul	0.03
9) Bis-(2-Chloroethyl)eth...	7.292	67	237466	30.762	ng/ul	0.00
11) 2-Chlorophenol-d4	7.504	132	278953	33.849	ng/ul	0.01
15) 4-Methylphenol-d8	8.716	113	222645	24.210	ng/ul	0.01
21) Nitrobenzene-d5	9.163	128	130835	30.855	ng/ul	0.01
24) 2-Nitrophenol-d4	9.898	143	162148	33.551	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.492	165	204568	24.443	ng/ul	0.05
31) 4-Chloroaniline-d4	11.069	131	43231	3.223	ng/ul	0.11
46) Dimethylphthalate-d6	14.033	166	1079941	35.060	ng/ul	0.00
49) Acenaphthylene-d8	14.310	160	1169283	34.638	ng/ul	0.00
54) 4-Nitrophenol-d4	15.039	143	1614	0.257	ng/ul	0.11
60) Fluorene-d10	15.621	176	956096	37.088	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.816	200	105012	19.639	ng/ul	0.02
73) Anthracene-d10	17.486	188	1648624	40.689	ng/ul	0.00
81) Pyrene-d10	19.715	212	2112449	40.559	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.797	264	2273049	40.607	ng/ul	0.00

Target Compounds Qvalue

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

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