

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100924\
 Data File : BM047954.D
 Acq On : 09 Oct 2024 21:21
 Operator : RC/JU
 Sample : P4187-16
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EY082

Quant Time: Oct 09 23:42:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 04 08:32:31 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.786	152	262017	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.580	136	1050433	20.000	ng/u1	#-0.01
38) Acenaphthene-d10	14.421	164	637397	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.162	188	1313611	20.000	ng/u1	-0.01
79) Chrysene-d12	21.386	240	1324445	20.000	ng/u1	-0.01
88) Perylene-d12	24.379	264	1586477	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	28043	3.459	ng/uL	-0.01
4) Pyridine-d5	3.663	84	125028	5.267	ng/u1	0.00
7) Phenol-d5	6.957	99	176857	7.138	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.116	67	366466	20.844	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.322	132	475588	26.383	ng/u1	-0.01
15) 4-Methylphenol-d8	8.498	113	319448	17.341	ng/u1	0.00
21) Nitrobenzene-d5	8.945	128	245779	29.387	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	268517	33.503	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.204	165	483103	30.174	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.716	131	158862	6.418	ng/u1	-0.01
46) Dimethylphthalate-d6	13.833	166	1360575	29.442	ng/u1	-0.01
49) Acenaphthylene-d8	14.115	160	1583334	28.898	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.627	143	60824	6.611	ng/u1	0.00
60) Fluorene-d10	15.415	176	1239682	30.971	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.533	200	217136	28.247	ng/u1	0.00
73) Anthracene-d10	17.256	188	2030888	31.442	ng/u1	-0.02
81) Pyrene-d10	19.550	212	2477906	32.965	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.174	264	2783467	33.815	ng/u1	-0.02

Target Compounds Qvalue

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

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