

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
 Data File : BM047091.D
 Acq On : 08 Aug 2024 18:06
 Operator : RC/JU
 Sample : P3435-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E05W8

Quant Time: Aug 08 18:54:24 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 07 04:48:25 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.387	152	218706	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.139	136	869417	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.045	164	546158	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.798	188	1129198	20.000	ng/u1	#-0.02
79) Chrysene-d12	21.033	240	992496	20.000	ng/u1	#-0.02
88) Perylene-d12	23.739	264	1200107	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.016	96	20514	3.716	ng/uL	0.00
4) Pyridine-d5	3.405	84	306645	20.364	ng/u1	0.00
7) Phenol-d5	6.593	99	414515	24.950	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	6.746	67	250317	22.898	ng/u1	-0.01
11) 2-Chlorophenol-d4	6.928	132	312796	22.226	ng/u1	-0.01
15) 4-Methylphenol-d8	8.104	113	329210	25.014	ng/u1	-0.02
21) Nitrobenzene-d5	8.534	128	147067	21.080	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.245	143	160477	21.063	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	9.781	165	314143	21.476	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.292	131	442238	23.004	ng/u1	-0.01
46) Dimethylphthalate-d6	13.469	166	996677	23.867	ng/u1	-0.01
49) Acenaphthylene-d8	13.727	160	1113926	23.897	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.286	143	130313	16.139	ng/u1	-0.01
60) Fluorene-d10	15.045	176	860043	24.075	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.186	200	88422	12.394	ng/u1	-0.01
73) Anthracene-d10	16.898	188	1386806	25.907	ng/u1	-0.01
81) Pyrene-d10	19.209	212	1522813	26.803	ng/u1	-0.02
92) Benzo(a)pyrene-d12	23.550	264	1642571	25.993	ng/u1	-0.02

Target Compounds Qvalue

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

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