

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
 Data File : BM047089.D
 Acq On : 08 Aug 2024 16:46
 Operator : RC/JU
 Sample : P3435-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E05W3

Quant Time: Aug 08 17:21:48 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	249793	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.139	136	998868	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.039	164	621789	20.000	ng/u1	-0.01
64) Phenanthrene-d10	16.804	188	1283276	20.000	ng/u1	-0.01
79) Chrysene-d12	21.033	240	1100696	20.000	ng/u1	#-0.02
88) Perylene-d12	23.738	264	1341804	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.016	96	18878	2.994	ng/uL	0.00
4) Pyridine-d5	3.404	84	276147	16.057	ng/u1	0.00
7) Phenol-d5	6.592	99	359281	18.934	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	6.745	67	226261	18.122	ng/u1	-0.01
11) 2-Chlorophenol-d4	6.928	132	278203	17.308	ng/u1	-0.01
15) 4-Methylphenol-d8	8.104	113	285350	18.983	ng/u1	-0.02
21) Nitrobenzene-d5	8.534	128	133664	16.676	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.245	143	145007	16.566	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	9.781	165	272814	16.233	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.292	131	394434	17.858	ng/u1	-0.01
46) Dimethylphthalate-d6	13.469	166	850483	17.889	ng/u1	-0.01
49) Acenaphthylene-d8	13.727	160	982140	18.507	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.286	143	107493	11.694	ng/u1	-0.01
60) Fluorene-d10	15.045	176	745585	18.333	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.186	200	74250	9.158	ng/u1	-0.01
73) Anthracene-d10	16.898	188	1176467	19.339	ng/u1	-0.01
81) Pyrene-d10	19.215	212	1290322	20.478	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.556	264	1359437	19.240	ng/u1	-0.02

Target Compounds Qvalue

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

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