

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
 Data File : BM047094.D
 Acq On : 08 Aug 2024 20:07
 Operator : RC/JU
 Sample : P3435-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E05X3

Quant Time: Aug 08 23:10:29 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	275879	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.139	136	1086565	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.039	164	685935	20.000	ng/u1	-0.01
64) Phenanthrene-d10	16.798	188	1384207	20.000	ng/u1	-0.02
79) Chrysene-d12	21.033	240	1162246	20.000	ng/u1	#-0.02
88) Perylene-d12	23.738	264	1320866	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.016	96	14317	2.056	ng/uL	0.00
4) Pyridine-d5	3.404	84	208910	10.999	ng/u1	0.00
7) Phenol-d5	6.593	99	274304	13.089	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	6.745	67	178416	12.939	ng/u1	-0.01
11) 2-Chlorophenol-d4	6.928	132	223282	12.578	ng/u1	-0.01
15) 4-Methylphenol-d8	8.104	113	214109	12.897	ng/u1	-0.02
21) Nitrobenzene-d5	8.534	128	103348	11.853	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.245	143	112048	11.768	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	9.781	165	215940	11.812	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.292	131	292187	12.161	ng/u1	-0.01
46) Dimethylphthalate-d6	13.469	166	637305	12.151	ng/u1	-0.01
49) Acenaphthylene-d8	13.727	160	747250	12.764	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.286	143	61926	6.107	ng/u1	-0.01
60) Fluorene-d10	15.045	176	562252	12.532	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.186	200	47305	5.409	ng/u1	-0.01
73) Anthracene-d10	16.898	188	845408	12.883	ng/u1	-0.01
81) Pyrene-d10	19.209	212	917238	13.786	ng/u1	-0.02
92) Benzo(a)pyrene-d12	23.550	264	895863	12.880	ng/u1	-0.02

Target Compounds Qvalue

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

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