

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
 Data File : BM047100.D
 Acq On : 09 Aug 2024 00:50
 Operator : RC/JU
 Sample : PB162448BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK448

Quant Time: Aug 09 02:06:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 09 01:30:34 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.387	152	225888	20.000	ng/u1	0.00
20) Naphthalene-d8	10.139	136	879328	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.039	164	544038	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.798	188	1093784	20.000	ng/u1	0.00
79) Chrysene-d12	21.033	240	905747	20.000	ng/u1	0.00
88) Perylene-d12	23.733	264	1016345	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.016	96	32671	5.730	ng/uL	0.00
4) Pyridine-d5	3.405	84	450646	28.976	ng/u1	0.00
7) Phenol-d5	6.593	99	526683	30.693	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.746	67	339238	30.046	ng/u1	0.00
11) 2-Chlorophenol-d4	6.928	132	412778	28.398	ng/u1	0.00
15) 4-Methylphenol-d8	8.104	113	401235	29.517	ng/u1	0.00
21) Nitrobenzene-d5	8.534	128	202923	28.759	ng/u1	0.00
24) 2-Nitrophenol-d4	9.245	143	217910	28.279	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.781	165	395522	26.735	ng/u1	0.00
31) 4-Chloroaniline-d4	10.292	131	545611	28.061	ng/u1	0.00
46) Dimethylphthalate-d6	13.469	166	1128389	27.126	ng/u1	0.00
49) Acenaphthylene-d8	13.727	160	1321987	28.471	ng/u1	0.00
54) 4-Nitrophenol-d4	14.280	143	191113	23.762	ng/u1	0.00
60) Fluorene-d10	15.045	176	967909	27.200	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.186	200	145822	21.102	ng/u1	0.00
73) Anthracene-d10	16.898	188	1495287	28.837	ng/u1	0.00
81) Pyrene-d10	19.209	212	1616920	31.185	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.550	264	1537134	28.722	ng/u1	0.00

Target Compounds Qvalue

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

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