

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
 Data File : BM046986.D
 Acq On : 05 Aug 2024 09:47
 Operator : RC/JU
 Sample : PB162449BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK449

Quant Time: Aug 06 06:25:38 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Aug 03 02:51:11 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.398	152	163068	20.000	ng/u1	0.00
20) Naphthalene-d8	10.151	136	613782	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.051	164	370353	20.000	ng/u1	-0.01
64) Phenanthrene-d10	16.815	188	751002	20.000	ng/u1	0.00
79) Chrysene-d12	21.050	240	662901	20.000	ng/u1	0.00
88) Perylene-d12	23.762	264	762777	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.016	96	25920	6.297	ng/uL	0.00
4) Pyridine-d5	3.405	84	343923	30.633	ng/u1	0.00
7) Phenol-d5	6.598	99	386422	31.195	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.751	67	257831	31.633	ng/u1	0.00
11) 2-Chlorophenol-d4	6.940	132	309910	29.534	ng/u1	0.00
15) 4-Methylphenol-d8	8.116	113	297539	30.321	ng/u1	0.00
21) Nitrobenzene-d5	8.539	128	153235	31.112	ng/u1	0.00
24) 2-Nitrophenol-d4	9.251	143	164761	30.633	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	9.786	165	296404	28.703	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.304	131	409789	30.194	ng/u1	0.00
46) Dimethylphthalate-d6	13.480	166	887129	31.328	ng/u1	0.00
49) Acenaphthylene-d8	13.739	160	988819	31.283	ng/u1	0.00
54) 4-Nitrophenol-d4	14.292	143	145192	26.518	ng/u1	0.00
60) Fluorene-d10	15.057	176	729187	30.102	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.192	200	125823	26.518	ng/u1	-0.01
73) Anthracene-d10	16.910	188	1121967	31.514	ng/u1	0.00
81) Pyrene-d10	19.227	212	1247178	32.866	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.580	264	1249519	31.109	ng/u1	0.00

Target Compounds Qvalue

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

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