

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
 Data File : BM045812.D
 Acq On : 04 Jun 2024 11:37
 Operator : MA/JU
 Sample : PB160908BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK908

Quant Time: Jun 05 00:01:01 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 31 13:01:11 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.469	152	341780	20.000	ng/u1	0.00
20) Naphthalene-d8	10.233	136	1469109	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.110	164	956704	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.851	188	1964611	20.000	ng/u1	0.00
79) Chrysene-d12	21.068	240	1579293	20.000	ng/u1	0.00
88) Perylene-d12	23.786	264	1523262	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.052	96	47302	5.366	ng/uL	0.00
4) Pyridine-d5	3.463	84	614610	27.077	ng/u1	0.00
7) Phenol-d5	6.716	99	879890	30.304	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.834	67	673400	32.298	ng/u1	0.00
11) 2-Chlorophenol-d4	7.028	132	700431	32.295	ng/u1	0.00
15) 4-Methylphenol-d8	8.234	113	697757	31.809	ng/u1	0.00
21) Nitrobenzene-d5	8.634	128	361793	31.497	ng/u1	0.00
24) 2-Nitrophenol-d4	9.345	143	369657	31.798	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.892	165	639287	29.325	ng/u1	0.00
31) 4-Chloroaniline-d4	10.416	131	968895	32.113	ng/u1	0.00
46) Dimethylphthalate-d6	13.557	166	2262762	34.164	ng/u1	0.00
49) Acenaphthylene-d8	13.798	160	2387513	31.043	ng/u1	0.00
54) 4-Nitrophenol-d4	14.404	143	331960	23.224	ng/u1	0.00
60) Fluorene-d10	15.110	176	1833765	32.454	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.245	200	227889	22.716	ng/u1	0.00
73) Anthracene-d10	16.957	188	2678926	31.367	ng/u1	0.00
81) Pyrene-d10	19.251	212	3144730	32.019	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.603	264	2269754	31.882	ng/u1	0.00

Target Compounds	Qvalue

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

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