

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081822\
 Data File : BM036348.D
 Acq On : 18 Aug 2022 16:29
 Operator : CG/JU
 Sample : PB146659BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK659

Quant Time: Aug 18 17:12:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 16 15:00:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.798	152	247089	20.000	ng/u1	0.00
20) Naphthalene-d8	10.598	136	1029099	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.439	164	553346	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.186	188	1010285	20.000	ng/u1	0.00
79) Chrysene-d12	21.368	240	913434	20.000	ng/u1	0.00
88) Perylene-d12	23.697	264	974618	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.204	96	30958	4.775	ng/uL	0.00
4) Pyridine-d5	3.634	84	353044	19.746	ng/u1	0.00
7) Phenol-d5	6.975	99	526349	24.956	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.134	67	349298	26.311	ng/u1	0.00
11) 2-Chlorophenol-d4	7.328	132	442008	26.673	ng/u1	0.00
15) 4-Methylphenol-d8	8.522	113	409659	26.014	ng/u1	0.00
21) Nitrobenzene-d5	8.969	128	213950	27.421	ng/u1	0.00
24) 2-Nitrophenol-d4	9.686	143	217430	27.473	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	363220	25.685	ng/u1	0.00
31) 4-Chloroaniline-d4	10.745	131	562168	25.643	ng/u1	-0.01
46) Dimethylphthalate-d6	13.863	166	1038540	29.667	ng/u1	0.00
49) Acenaphthylene-d8	14.133	160	1300473	29.785	ng/u1	0.00
54) 4-Nitrophenol-d4	14.674	143	139071	20.966	ng/u1	0.00
60) Fluorene-d10	15.433	176	901396	28.495	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	110116	21.201	ng/u1	0.00
73) Anthracene-d10	17.280	188	1343564	29.909	ng/u1	0.00
81) Pyrene-d10	19.574	212	1454760	30.702	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.544	264	1439034	30.115	ng/u1	0.00

Target Compounds Qvalue

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

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