

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072622\
 Data File : BM035918.D
 Acq On : 26 Jul 2022 10:17
 Operator : CG/JU
 Sample : PB146421BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK421

Quant Time: Jul 26 14:08:29 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 25 14:55:39 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	432067	20.000	ng/u1	0.00
20) Naphthalene-d8	10.692	136	1913616	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.521	164	1202630	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	2505705	20.000	ng/u1	0.00
79) Chrysene-d12	21.433	240	2303969	20.000	ng/u1	0.00
88) Perylene-d12	23.809	264	2170605	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	80081	6.974	ng/uL	0.00
4) Pyridine-d5	3.699	84	1098828	34.932	ng/u1	0.00
7) Phenol-d5	7.045	99	1245408	34.372	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	852290	35.285	ng/u1	0.00
11) 2-Chlorophenol-d4	7.416	132	1076853	35.805	ng/u1	0.00
15) 4-Methylphenol-d8	8.598	113	1017412	34.661	ng/u1	0.00
21) Nitrobenzene-d5	9.051	128	553822	36.523	ng/u1	0.00
24) 2-Nitrophenol-d4	9.775	143	552673	37.410	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	926534	34.676	ng/u1	0.00
31) 4-Chloroaniline-d4	10.828	131	1594842	35.258	ng/u1	0.00
46) Dimethylphthalate-d6	13.939	166	3285084	40.101	ng/u1	0.00
49) Acenaphthylene-d8	14.216	160	4034380	38.598	ng/u1	0.00
54) 4-Nitrophenol-d4	14.716	143	582641	34.635	ng/u1	0.00
60) Fluorene-d10	15.510	176	2861820	38.659	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.627	200	434720	32.276	ng/u1	0.00
73) Anthracene-d10	17.357	188	4372040	38.593	ng/u1	0.00
81) Pyrene-d10	19.651	212	4983043	39.024	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.656	264	4296653	38.990	ng/u1	0.00

Target Compounds Qvalue

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

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