

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072922\
 Data File : BM035995.D
 Acq On : 29 Jul 2022 16:40
 Operator : CG/JU
 Sample : PB146498BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK498

Quant Time: Jul 30 00:24:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 30 00:20:59 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	279378	20.000	ng/u1	0.00
20) Naphthalene-d8	10.680	136	1158328	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.515	164	673298	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.251	188	1369422	20.000	ng/u1	0.00
79) Chrysene-d12	21.433	240	1287904	20.000	ng/u1	0.00
88) Perylene-d12	23.797	264	1317445	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	41159	5.544	ng/uL	0.00
4) Pyridine-d5	3.693	84	529053	26.011	ng/u1	0.00
7) Phenol-d5	7.039	99	627978	26.804	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	404299	25.886	ng/u1	0.00
11) 2-Chlorophenol-d4	7.404	132	555601	28.570	ng/u1	0.00
15) 4-Methylphenol-d8	8.586	113	492761	25.962	ng/u1	0.00
21) Nitrobenzene-d5	9.045	128	259730	28.297	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	269719	30.161	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.304	165	455251	28.148	ng/u1	0.00
31) 4-Chloroaniline-d4	10.822	131	717201	26.194	ng/u1	0.00
46) Dimethylphthalate-d6	13.927	166	1353921	29.521	ng/u1	0.00
49) Acenaphthylene-d8	14.210	160	1723187	29.448	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	223379	23.718	ng/u1	0.00
60) Fluorene-d10	15.504	176	1193991	28.810	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.621	200	160542	21.810	ng/u1	0.00
73) Anthracene-d10	17.351	188	1827476	29.517	ng/u1	0.00
81) Pyrene-d10	19.639	212	2126429	29.791	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.650	264	2026597	30.300	ng/u1	0.00

Target Compounds Qvalue

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

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