

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072622\
 Data File : BM035930.D
 Acq On : 26 Jul 2022 17:42
 Operator : CG/JU
 Sample : PB146378BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK378

Quant Time: Jul 26 18:48:13 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	289020	20.000	ng/u1	0.00
20) Naphthalene-d8	10.692	136	1314961	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.521	164	843296	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	1793270	20.000	ng/u1	0.00
79) Chrysene-d12	21.433	240	1767750	20.000	ng/u1	0.00
88) Perylene-d12	23.803	264	1837123	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	52675	6.858	ng/uL	0.00
4) Pyridine-d5	3.693	84	737222	35.036	ng/u1	0.00
7) Phenol-d5	7.045	99	843391	34.798	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	591974	36.638	ng/u1	0.00
11) 2-Chlorophenol-d4	7.416	132	712867	35.434	ng/u1	0.00
15) 4-Methylphenol-d8	8.598	113	689053	35.093	ng/u1	0.00
21) Nitrobenzene-d5	9.051	128	364496	34.981	ng/u1	0.00
24) 2-Nitrophenol-d4	9.775	143	355425	35.011	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	626837	34.140	ng/u1	0.00
31) 4-Chloroaniline-d4	10.833	131	1095508	35.245	ng/u1	0.00
46) Dimethylphthalate-d6	13.939	166	2287686	39.825	ng/u1	0.00
49) Acenaphthylene-d8	14.216	160	2747540	37.488	ng/u1	0.00
54) 4-Nitrophenol-d4	14.716	143	396884	33.646	ng/u1	0.00
60) Fluorene-d10	15.510	176	2013677	38.793	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.627	200	273676	28.392	ng/u1	0.00
73) Anthracene-d10	17.357	188	3105600	38.305	ng/u1	0.00
81) Pyrene-d10	19.645	212	3668319	37.442	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.656	264	3559334	38.162	ng/u1	0.00

Target Compounds Qvalue

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

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