

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072622\
 Data File : BM035929.D
 Acq On : 26 Jul 2022 17:05
 Operator : CG/JU
 Sample : PB146343BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK343

Quant Time: Jul 26 17:38:01 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	311657	20.000	ng/u1	0.00
20) Naphthalene-d8	10.693	136	1392479	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.522	164	893013	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	1892813	20.000	ng/u1	0.00
79) Chrysene-d12	21.433	240	1834770	20.000	ng/u1	0.00
88) Perylene-d12	23.809	264	1859178	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	56486	6.820	ng/uL	0.00
4) Pyridine-d5	3.693	84	771266	33.992	ng/u1	0.00
7) Phenol-d5	7.046	99	888573	33.999	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	616008	35.356	ng/u1	0.00
11) 2-Chlorophenol-d4	7.410	132	750914	34.614	ng/u1	0.00
15) 4-Methylphenol-d8	8.599	113	717104	33.869	ng/u1	0.00
21) Nitrobenzene-d5	9.052	128	383735	34.777	ng/u1	0.00
24) 2-Nitrophenol-d4	9.775	143	373355	34.730	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	649226	33.391	ng/u1	0.00
31) 4-Chloroaniline-d4	10.828	131	1146918	34.845	ng/u1	0.00
46) Dimethylphthalate-d6	13.939	166	2386029	39.224	ng/u1	0.00
49) Acenaphthylene-d8	14.216	160	2879517	37.101	ng/u1	0.00
54) 4-Nitrophenol-d4	14.716	143	413584	33.109	ng/u1	0.00
60) Fluorene-d10	15.510	176	2087737	37.981	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	288595	28.365	ng/u1	0.00
73) Anthracene-d10	17.357	188	3218025	37.605	ng/u1	0.00
81) Pyrene-d10	19.651	212	3759191	36.968	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.656	264	3574368	37.869	ng/u1	0.00

Target Compounds Qvalue

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

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