

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
 Data File : BM034159.D
 Acq On : 08 Mar 2022 21:41
 Operator : CG/JU
 Sample : N1733-08
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBRS3

Quant Time: Mar 09 00:42:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 08 13:09:16 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.989	152	156088	20.000	ng/u1	0.00
20) Naphthalene-d8	10.807	136	620063	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.630	164	426735	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.365	188	880116	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.536	240	771850	20.000	ng/u1	# 0.00
88) Perylene-d12	23.983	264	801616	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	27900	8.221	ng/uL	0.00
4) Pyridine-d5	3.790	84	379187	38.383	ng/u1	0.00
7) Phenol-d5	7.142	99	495782	40.537	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.313	67	281103	42.702	ng/u1	0.00
11) 2-Chlorophenol-d4	7.519	132	437807	43.184	ng/u1	0.00
15) 4-Methylphenol-d8	8.695	113	427163	41.683	ng/u1	0.00
21) Nitrobenzene-d5	9.154	128	217878	45.350	ng/u1	0.00
24) 2-Nitrophenol-d4	9.878	143	250387	47.985	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.419	165	462148	42.947	ng/u1	0.00
31) 4-Chloroaniline-d4	10.936	131	600972	40.876	ng/u1	0.00
46) Dimethylphthalate-d6	14.036	166	1439370	44.449	ng/u1	0.00
49) Acenaphthylene-d8	14.324	160	1777109	43.886	ng/u1	0.00
54) 4-Nitrophenol-d4	14.807	143	215540	35.772	ng/u1	0.00
60) Fluorene-d10	15.618	176	1249110	43.342	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.724	200	212068	37.338	ng/u1	0.00
73) Anthracene-d10	17.465	188	1920200	44.291	ng/u1	0.00
81) Pyrene-d10	19.753	212	2069038	48.768	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.830	264	1886821	44.371	ng/u1	0.00

Target Compounds Qvalue

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

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