

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072622\
 Data File : BM035932.D
 Acq On : 26 Jul 2022 18:56
 Operator : CG/JU
 Sample : N3792-17
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Jul 26 23:42:17 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	384436	20.000	ng/u1	0.00
20) Naphthalene-d8	10.692	136	1764442	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.522	164	1137349	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	2464899	20.000	ng/u1	0.00
79) Chrysene-d12	21.433	240	2399388	20.000	ng/u1	0.00
88) Perylene-d12	23.803	264	2197990	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	44815	4.387	ng/uL	0.00
4) Pyridine-d5	3.705	84	134229	4.796	ng/u1	0.00
7) Phenol-d5	7.046	99	251983	7.816	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	753682	35.069	ng/u1	0.00
11) 2-Chlorophenol-d4	7.410	132	720014	26.907	ng/u1	0.00
15) 4-Methylphenol-d8	8.593	113	459194	17.582	ng/u1	0.00
21) Nitrobenzene-d5	9.051	128	482173	34.486	ng/u1	0.00
24) 2-Nitrophenol-d4	9.775	143	431598	31.684	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	757094	30.730	ng/u1	0.00
31) 4-Chloroaniline-d4	10.828	131	549177	13.167	ng/u1	0.00
46) Dimethylphthalate-d6	13.939	166	2937288	37.913	ng/u1	0.00
49) Acenaphthylene-d8	14.216	160	3496232	35.370	ng/u1	0.00
54) 4-Nitrophenol-d4	14.716	143	92273	5.800	ng/u1	0.00
60) Fluorene-d10	15.510	176	2609144	37.269	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.621	200	154103	11.631	ng/u1	0.00
73) Anthracene-d10	17.357	188	4040746	36.260	ng/u1	0.00
81) Pyrene-d10	19.645	212	5049945	37.975	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.656	264	4263968	38.211	ng/u1	0.00

Target Compounds	Qvalue

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

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