

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
 Data File : BM035951.D
 Acq On : 27 Jul 2022 15:21
 Operator : CG/JU
 Sample : N3795-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Jul 27 23:12:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 27 03:28:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	259090	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.698	136	1080384	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.527	164	654513	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.268	188	1335076	20.000	ng/u1	-0.01
79) Chrysene-d12	21.445	240	1218784	20.000	ng/u1	-0.02
88) Perylene-d12	23.827	264	1297634	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	34017	4.941	ng/uL	0.00
4) Pyridine-d5	3.699	84	448049	23.753	ng/u1	0.00
7) Phenol-d5	7.057	99	640559	29.482	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.222	67	400963	27.683	ng/u1	0.00
11) 2-Chlorophenol-d4	7.416	132	550461	30.522	ng/u1	-0.01
15) 4-Methylphenol-d8	8.604	113	557457	31.671	ng/u1	-0.02
21) Nitrobenzene-d5	9.057	128	287778	33.615	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.781	143	310633	37.243	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.322	165	527581	34.973	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.839	131	786495	30.797	ng/u1	-0.01
46) Dimethylphthalate-d6	13.945	166	1869693	41.936	ng/u1	0.00
49) Acenaphthylene-d8	14.221	160	2302138	40.470	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.733	143	333450	36.421	ng/u1	-0.02
60) Fluorene-d10	15.521	176	1625187	40.339	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	241065	33.591	ng/u1	-0.02
73) Anthracene-d10	17.368	188	2523726	41.812	ng/u1	-0.01
81) Pyrene-d10	19.656	212	2892563	42.823	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.674	264	2769051	42.032	ng/u1	-0.02

Target Compounds	Qvalue

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

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