

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
 Data File : BM035952.D
 Acq On : 27 Jul 2022 15:58
 Operator : CG/JU
 Sample : N3741-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H0B12

Quant Time: Jul 27 23:13:07 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	270925	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.698	136	1134901	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.527	164	669660	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.268	188	1355646	20.000	ng/u1	-0.01
79) Chrysene-d12	21.445	240	1176727	20.000	ng/u1	-0.02
88) Perylene-d12	23.827	264	1231010	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	31055	4.313	ng/uL	0.00
4) Pyridine-d5	3.705	84	128814	6.531	ng/u1	0.00
7) Phenol-d5	7.057	99	544775	23.978	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.222	67	334383	22.078	ng/u1	0.00
11) 2-Chlorophenol-d4	7.416	132	475693	25.224	ng/u1	-0.01
15) 4-Methylphenol-d8	8.604	113	407925	22.163	ng/u1	-0.02
21) Nitrobenzene-d5	9.057	128	230554	25.637	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.781	143	240914	27.496	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.316	165	408596	25.784	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.839	131	611478	22.794	ng/u1	-0.01
46) Dimethylphthalate-d6	13.939	166	1300990	28.521	ng/u1	-0.01
49) Acenaphthylene-d8	14.222	160	1596781	27.436	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.727	143	165867	17.707	ng/u1	-0.02
60) Fluorene-d10	15.516	176	1136041	27.560	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.645	200	113155	15.528	ng/u1	-0.02
73) Anthracene-d10	17.368	188	1722238	28.100	ng/u1	-0.01
81) Pyrene-d10	19.656	212	1981645	30.386	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.674	264	1828004	29.250	ng/u1	-0.02

Target Compounds Qvalue

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

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