

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072722\
 Data File : BM035958.D
 Acq On : 27 Jul 2022 19:39
 Operator : CG/JU
 Sample : N3874-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Jul 27 23:14:31 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	7.886	152	253322	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.692	136	1063199	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.521	164	630431	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.262	188	1289492	20.000	ng/u1	-0.02
79) Chrysene-d12	21.439	240	1158791	20.000	ng/u1	-0.02
88) Perylene-d12	23.815	264	1207717	20.000	ng/u1	-0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	33326	4.950	ng/uL	0.00
4) Pyridine-d5	3.710	84	70970	3.848	ng/u1	0.00
7) Phenol-d5	7.051	99	178279	8.392	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.222	67	435485	30.751	ng/u1	0.00
11) 2-Chlorophenol-d4	7.416	132	488704	27.715	ng/u1	-0.01
15) 4-Methylphenol-d8	8.598	113	317779	18.465	ng/u1	-0.02
21) Nitrobenzene-d5	9.051	128	304813	36.180	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.775	143	307409	37.452	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.316	165	486008	32.738	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.833	131	623236	24.799	ng/u1	-0.02
46) Dimethylphthalate-d6	13.939	166	1729875	40.283	ng/u1	-0.01
49) Acenaphthylene-d8	14.215	160	2110282	38.515	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.721	143	74231	8.418	ng/u1	-0.03
60) Fluorene-d10	15.515	176	1497796	38.597	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.639	200	243553	35.138	ng/u1	-0.03
73) Anthracene-d10	17.362	188	2381343	40.847	ng/u1	-0.02
81) Pyrene-d10	19.650	212	2714401	42.266	ng/u1	-0.02
92) Benzo(a)pyrene-d12	23.668	264	2517576	41.060	ng/u1	-0.02

Target Compounds Qvalue

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

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