

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072922\
 Data File : BM036006.D
 Acq On : 29 Jul 2022 23:23
 Operator : CG/JU
 Sample : N3790-19
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBUL5

Quant Time: Jul 30 00:25:19 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 30 00:20:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	275441	20.000	ng/u1	0.00
20) Naphthalene-d8	10.680	136	1162878	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.510	164	660139	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.251	188	1271014	20.000	ng/u1	0.00
79) Chrysene-d12	21.427	240	1155484	20.000	ng/u1	0.00
88) Perylene-d12	23.791	264	1162716	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	41451	5.663	ng/uL	0.00
4) Pyridine-d5	3.693	84	485705	24.221	ng/u1	0.00
7) Phenol-d5	7.040	99	732861	31.728	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	466733	30.311	ng/u1	0.00
11) 2-Chlorophenol-d4	7.404	132	620279	32.352	ng/u1	0.00
15) 4-Methylphenol-d8	8.586	113	570023	30.462	ng/u1	0.00
21) Nitrobenzene-d5	9.039	128	300698	32.632	ng/u1	0.00
24) 2-Nitrophenol-d4	9.763	143	310411	34.576	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.298	165	539370	33.218	ng/u1	0.00
31) 4-Chloroaniline-d4	10.822	131	814368	29.627	ng/u1	0.00
46) Dimethylphthalate-d6	13.927	166	1584126	35.228	ng/u1	0.00
49) Acenaphthylene-d8	14.204	160	2016476	35.146	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	231171	25.035	ng/u1	0.00
60) Fluorene-d10	15.504	176	1400453	34.465	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.621	200	152881	22.377	ng/u1	0.00
73) Anthracene-d10	17.351	188	2049143	35.660	ng/u1	0.00
81) Pyrene-d10	19.639	212	2233735	34.881	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.638	264	2104997	35.660	ng/u1	-0.01

Target Compounds	Qvalue

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

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