

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072922\
 Data File : BM036022.D
 Acq On : 30 Jul 2022 09:38
 Operator : CG/JU
 Sample : N3792-14
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Aug 01 00:46:07 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	7.869	152	319719	20.000	ng/u1	0.00
20) Naphthalene-d8	10.675	136	1317259	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.510	164	734610	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.251	188	1448602	20.000	ng/u1	0.00
79) Chrysene-d12	21.427	240	1420729	20.000	ng/u1	0.00
88) Perylene-d12	23.791	264	1463495	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	17275	2.033	ng/uL	0.00
4) Pyridine-d5	3.699	84	200477	8.613	ng/u1	0.00
7) Phenol-d5	7.034	99	350754	13.082	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.204	67	686574	38.413	ng/u1	0.00
11) 2-Chlorophenol-d4	7.398	132	355835	15.989	ng/u1	-0.01
15) 4-Methylphenol-d8	8.581	113	402121	18.513	ng/u1	-0.01
21) Nitrobenzene-d5	9.039	128	442225	42.366	ng/u1	0.00
24) 2-Nitrophenol-d4	9.757	143	124539	12.246	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.298	165	269220	14.637	ng/u1	0.00
31) 4-Chloroaniline-d4	10.816	131	962951	30.926	ng/u1	-0.01
46) Dimethylphthalate-d6	13.927	166	2145221	42.870	ng/u1	0.00
49) Acenaphthylene-d8	14.204	160	2765447	43.314	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	43797	4.262	ng/u1	0.00
60) Fluorene-d10	15.498	176	1966679	43.493	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.615	200	18060	2.319	ng/u1	-0.01
73) Anthracene-d10	17.345	188	3073986	46.937	ng/u1	0.00
81) Pyrene-d10	19.639	212	3555481	45.155	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.644	264	3557306	47.878	ng/u1	0.00

Target Compounds	Qvalue

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

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