

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110222\
 Data File : BM037322.D
 Acq On : 02 Nov 2022 18:41
 Operator : CG/JU
 Sample : N5301-11
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBWX6

Quant Time: Nov 02 22:55:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 02 03:33:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	306527	20.000	ng/u1	0.00
20) Naphthalene-d8	10.657	136	1378487	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.492	164	880692	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1838442	20.000	ng/u1	0.00
79) Chrysene-d12	21.403	240	1820955	20.000	ng/u1	0.00
88) Perylene-d12	23.750	264	1693910	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	31141	4.478	ng/uL	0.00
4) Pyridine-d5	3.699	84	212479	10.236	ng/u1	0.00
7) Phenol-d5	7.028	99	184648	6.922	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.192	67	575675	35.167	ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	543631	26.672	ng/u1	0.00
15) 4-Methylphenol-d8	8.575	113	361061	16.575	ng/u1	0.00
21) Nitrobenzene-d5	9.028	128	362875	35.240	ng/u1	0.00
24) 2-Nitrophenol-d4	9.745	143	367107	32.729	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.280	165	625707	30.258	ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	695962	22.023	ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	2241413	36.809	ng/u1	0.00
49) Acenaphthylene-d8	14.186	160	2670675	36.481	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	52702	4.330	ng/u1	0.00
60) Fluorene-d10	15.480	176	2121455	39.068	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.610	200	313623	26.934	ng/u1	0.00
73) Anthracene-d10	17.327	188	3389173	40.029	ng/u1	0.00
81) Pyrene-d10	19.621	212	3874065	40.713	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.603	264	3469172	39.786	ng/u1	0.00

Target Compounds Qvalue

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

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