

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110222\
 Data File : BM037313.D
 Acq On : 02 Nov 2022 13:13
 Operator : CG/JU
 Sample : N5301-05
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBWP1

Quant Time: Nov 02 22:53:02 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	7.845	152	354724	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.651	136	1527473	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.486	164	934782	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1829045	20.000	ng/u1	0.00
79) Chrysene-d12	21.403	240	1578990	20.000	ng/u1	0.00
88) Perylene-d12	23.750	264	1525453	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	33735	4.192	ng/uL	0.00
4) Pyridine-d5	3.699	84	225372	9.382	ng/u1	0.00
7) Phenol-d5	7.022	99	174478	5.652	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	538977	28.452	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	507964	21.536	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	325423	12.909	ng/u1	-0.01
21) Nitrobenzene-d5	9.016	128	322143	28.233	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.739	143	329943	26.547	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	558397	24.369	ng/u1	0.00
31) 4-Chloroaniline-d4	10.798	131	784474	22.403	ng/u1	0.00
46) Dimethylphthalate-d6	13.904	166	1924132	29.770	ng/u1	0.00
49) Acenaphthylene-d8	14.180	160	2260417	29.090	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	49784	3.854	ng/u1	0.00
60) Fluorene-d10	15.480	176	1757003	30.484	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	259316	22.384	ng/u1	0.00
73) Anthracene-d10	17.327	188	2711171	32.185	ng/u1	0.00
81) Pyrene-d10	19.615	212	2859994	34.662	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.597	264	2535821	32.294	ng/u1	0.00

Target Compounds	Qvalue

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

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