

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
 Data File : BM037349.D
 Acq On : 03 Nov 2022 16:39
 Operator : CG/JU
 Sample : N5276-08
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4W6

Quant Time: Nov 04 01:04:09 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	7.851	152	293592	20.000	ng/u1	0.00
20) Naphthalene-d8	10.651	136	1326509	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.486	164	863810	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1802924	20.000	ng/u1	0.00
79) Chrysene-d12	21.397	240	1600322	20.000	ng/u1	0.00
88) Perylene-d12	23.750	264	1332760	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	27285	4.097	ng/uL	0.00
4) Pyridine-d5	3.704	84	498	0.025	ng/u1	0.00
7) Phenol-d5	7.028	99	169182	6.621	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.192	67	503857	32.136	ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	495827	25.399	ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	326992	15.672	ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	325316	32.831	ng/u1	0.00
24) 2-Nitrophenol-d4	9.739	143	336482	31.174	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.280	165	570656	28.677	ng/u1	0.00
31) 4-Chloroaniline-d4	10.816	131	35577	1.170	ng/u1	0.01
46) Dimethylphthalate-d6	13.910	166	2098944	35.143	ng/u1	0.00
49) Acenaphthylene-d8	14.186	160	2410433	33.570	ng/u1	0.00
54) 4-Nitrophenol-d4	14.715	143	58345	4.887	ng/u1	0.00
60) Fluorene-d10	15.480	176	1964144	36.878	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	319846	28.009	ng/u1	0.00
73) Anthracene-d10	17.327	188	3239462	39.014	ng/u1	0.00
81) Pyrene-d10	19.615	212	3554421	42.504	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.597	264	2706766	39.455	ng/u1	0.00

Target Compounds Qvalue

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

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