

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
 Data File : BM037370.D
 Acq On : 04 Nov 2022 05:40
 Operator : CG/JU
 Sample : N5276-18
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4X8

Quant Time: Nov 04 06:27:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 04 02:39:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.845	152	317296	20.000	ng/u1	0.00
20) Naphthalene-d8	10.645	136	1447084	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.480	164	944043	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.221	188	1908542	20.000	ng/u1	0.00
79) Chrysene-d12	21.392	240	1699740	20.000	ng/u1	0.00
88) Perylene-d12	23.738	264	1433690	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	35690	4.958	ng/uL	0.00
4) Pyridine-d5	3.699	84	180694	8.410	ng/u1	0.00
7) Phenol-d5	7.022	99	247194	8.952	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	576993	34.051	ng/u1	0.00
11) 2-Chlorophenol-d4	7.375	132	610455	28.934	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	422418	18.733	ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	384807	35.599	ng/u1	0.00
24) 2-Nitrophenol-d4	9.733	143	409268	34.758	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.269	165	723354	33.321	ng/u1	0.00
31) 4-Chloroaniline-d4	10.792	131	460736	13.888	ng/u1	0.00
46) Dimethylphthalate-d6	13.898	166	2497652	38.265	ng/u1	0.00
49) Acenaphthylene-d8	14.174	160	2982028	38.001	ng/u1	0.00
54) 4-Nitrophenol-d4	14.704	143	104627	8.019	ng/u1	0.00
60) Fluorene-d10	15.474	176	2317361	39.812	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	378182	31.285	ng/u1	0.00
73) Anthracene-d10	17.321	188	3504900	39.875	ng/u1	0.00
81) Pyrene-d10	19.609	212	4010620	45.154	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.586	264	2947733	39.942	ng/u1	0.00

Target Compounds Qvalue

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

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