

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110222\
 Data File : BM037317.D
 Acq On : 02 Nov 2022 15:39
 Operator : CG/JU
 Sample : N5301-10
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBWU8

Quant Time: Nov 02 22:53:39 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	324978	20.000	ng/u1	0.00
20) Naphthalene-d8	10.657	136	1334562	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.492	164	770200	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	1423707	20.000	ng/u1	0.00
79) Chrysene-d12	21.403	240	1309480	20.000	ng/u1	0.00
88) Perylene-d12	23.750	264	1373247	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	32832	4.454	ng/uL	0.00
4) Pyridine-d5	3.699	84	192557	8.750	ng/u1	0.00
7) Phenol-d5	7.028	99	186282	6.587	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	575483	33.159	ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	566084	26.197	ng/u1	0.00
15) 4-Methylphenol-d8	8.575	113	343752	14.884	ng/u1	0.00
21) Nitrobenzene-d5	9.028	128	351377	35.247	ng/u1	0.00
24) 2-Nitrophenol-d4	9.745	143	357088	32.884	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	592871	29.613	ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	747281	24.425	ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	1801229	33.824	ng/u1	0.00
49) Acenaphthylene-d8	14.186	160	2225344	34.759	ng/u1	0.00
54) 4-Nitrophenol-d4	14.716	143	42429	3.986	ng/u1	0.00
60) Fluorene-d10	15.480	176	1687265	35.529	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.610	200	245437	27.218	ng/u1	0.00
73) Anthracene-d10	17.327	188	2475270	37.751	ng/u1	0.00
81) Pyrene-d10	19.621	212	2659565	38.867	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.603	264	2675870	37.854	ng/u1	0.00

Target Compounds Qvalue

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

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