

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
 Data File : BM036658.D
 Acq On : 22 Sep 2022 12:14
 Operator : CG/JU
 Sample : N4708-03
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB8J5

Quant Time: Sep 23 00:59:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 21 22:42:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.022	152	377970	20.000	ng/u1	0.00
20) Naphthalene-d8	10.833	136	1704449	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.645	164	1070935	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.380	188	2113931	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	1520730	20.000	ng/u1	0.00
88) Perylene-d12	23.986	264	1333115	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	41841	4.150	ng/uL	0.00
4) Pyridine-d5	3.828	84	213368	7.256	ng/u1	0.01
7) Phenol-d5	7.175	99	213230	6.313	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.345	67	540396	26.344	ng/u1	0.00
11) 2-Chlorophenol-d4	7.545	132	552840	22.734	ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	379982	15.225	ng/u1	0.00
21) Nitrobenzene-d5	9.186	128	351841	30.498	ng/u1	0.00
24) 2-Nitrophenol-d4	9.910	143	329634	31.899	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	622776	26.005	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	519411	13.045	ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	2197928	30.430	ng/u1	0.00
49) Acenaphthylene-d8	14.339	160	2485735	29.013	ng/u1	0.00
54) 4-Nitrophenol-d4	14.827	143	75139	5.363	ng/u1	0.00
60) Fluorene-d10	15.633	176	1954293	30.192	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	233538	27.923	ng/u1	0.00
73) Anthracene-d10	17.480	188	3198284	33.090	ng/u1	0.00
81) Pyrene-d10	19.762	212	3481975	46.542	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.833	264	2421896	37.185	ng/u1	0.00

Target Compounds	Qvalue

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

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