

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122022\
 Data File : BM038171.D
 Acq On : 21 Dec 2022 14:35
 Operator : CG/JU
 Sample : N6014-17
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Dec 21 23:32:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 21 00:11:56 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.039	152	193402	20.000	ng/ul	0.00
20) Naphthalene-d8	10.857	136	798719	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.668	164	484801	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.409	188	971840	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.574	240	686555	20.000	ng/ul	0.00
88) Perylene-d12	24.027	264	726228	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	16231	3.811	ng/uL	0.00
4) Pyridine-d5	3.834	84	224881	17.797	ng/ul	0.00
7) Phenol-d5	7.204	99	334119	21.260	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.369	67	189617	19.814	ng/ul	0.00
11) 2-Chlorophenol-d4	7.569	132	303814	23.565	ng/ul	0.00
15) 4-Methylphenol-d8	8.757	113	269276	21.950	ng/ul	0.00
21) Nitrobenzene-d5	9.216	128	151222	23.913	ng/ul	0.00
24) 2-Nitrophenol-d4	9.939	143	174828	26.380	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.475	165	324443	25.775	ng/ul	0.00
31) 4-Chloroaniline-d4	10.998	131	390251	22.460	ng/ul	0.00
46) Dimethylphthalate-d6	14.080	166	940960	27.224	ng/ul	0.00
49) Acenaphthylene-d8	14.368	160	1081761	25.608	ng/ul	0.00
54) 4-Nitrophenol-d4	14.868	143	134348	22.808	ng/ul	0.00
60) Fluorene-d10	15.657	176	802557	26.063	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.780	200	70166	13.504	ng/ul	0.00
73) Anthracene-d10	17.509	188	1183094	26.084	ng/ul	0.00
81) Pyrene-d10	19.792	212	1171360	28.403	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.868	264	955421	26.339	ng/ul	0.00

Target Compounds Qvalue

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

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