

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
 Data File : BM036694.D
 Acq On : 24 Sep 2022 01:59
 Operator : CG/JU
 Sample : N4649-12
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EXZ17

Quant Time: Sep 24 02:54:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 24 00:41:16 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	445508	20.000	ng/ul	0.00
20) Naphthalene-d8	10.822	136	1999395	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.633	164	1217987	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.374	188	2226022	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	1551754	20.000	ng/ul	0.00
88) Perylene-d12	23.974	264	1401219	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	45849	3.858	ng/uL	0.00
4) Pyridine-d5	3.822	84	255043	7.359	ng/ul	0.00
7) Phenol-d5	7.169	99	242026	6.079	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	689199	28.505	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	691839	24.137	ng/ul	0.00
15) 4-Methylphenol-d8	8.716	113	444979	15.126	ng/ul	-0.01
21) Nitrobenzene-d5	9.175	128	440531	32.553	ng/ul	0.00
24) 2-Nitrophenol-d4	9.898	143	410452	33.861	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.439	165	736564	26.220	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	1117181	23.919	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166	2577655	31.379	ng/ul	0.00
49) Acenaphthylene-d8	14.333	160	2957198	30.349	ng/ul	0.00
54) 4-Nitrophenol-d4	14.821	143	61963	3.889	ng/ul	-0.01
60) Fluorene-d10	15.621	176	2299080	31.231	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.733	200	198401	22.527	ng/ul	0.00
73) Anthracene-d10	17.468	188	3501450	34.402	ng/ul	0.00
81) Pyrene-d10	19.751	212	3410101	44.670	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	2516890	36.766	ng/ul	0.00

Target Compounds Qvalue

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

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