

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
 Data File : BM036697.D
 Acq On : 24 Sep 2022 03:45
 Operator : CG/JU
 Sample : N4649-18
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EXZ22

Quant Time: Sep 24 04:35:44 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	521056	20.000	ng/ul	0.00
20) Naphthalene-d8	10.822	136	2306847	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.633	164	1403351	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.368	188	2573867	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	1789373	20.000	ng/ul	-0.01
88) Perylene-d12	23.974	264	1704646	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	55406	3.986	ng/uL	0.00
4) Pyridine-d5	3.816	84	252021	6.217	ng/ul	0.00
7) Phenol-d5	7.169	99	223277	4.795	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	603397	21.338	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	626088	18.676	ng/ul	0.00
15) 4-Methylphenol-d8	8.716	113	407837	11.854	ng/ul	-0.01
21) Nitrobenzene-d5	9.175	128	381066	24.406	ng/ul	0.00
24) 2-Nitrophenol-d4	9.898	143	354347	25.336	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	659304	20.341	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	817132	15.163	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	2165459	22.879	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	2557742	22.782	ng/ul	0.00
54) 4-Nitrophenol-d4	14.822	143	56100	3.056	ng/ul	-0.01
60) Fluorene-d10	15.622	176	1949389	22.983	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.733	200	170698	16.762	ng/ul	0.00
73) Anthracene-d10	17.468	188	2968264	25.222	ng/ul	0.00
81) Pyrene-d10	19.751	212	3012541	34.222	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.815	264	2341354	28.114	ng/ul	0.00

Target Compounds Qvalue

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

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