

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
 Data File : BM036727.D
 Acq On : 24 Sep 2022 23:00
 Operator : CG/JU
 Sample : N4662-01
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EXZ26

Quant Time: Sep 26 01:15:18 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	8.004	152	377366	20.000	ng/u1	0.00
20) Naphthalene-d8	10.816	136	1703375	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.627	164	1058937	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.368	188	2013051	20.000	ng/u1	0.00
79) Chrysene-d12	21.533	240	1364019	20.000	ng/u1	-0.01
88) Perylene-d12	23.962	264	1205169	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	48880	4.856	ng/uL	0.00
4) Pyridine-d5	3.816	84	200769	6.839	ng/u1	0.00
7) Phenol-d5	7.163	99	219408	6.506	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.334	67	565134	27.595	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	583782	24.045	ng/u1	0.00
15) 4-Methylphenol-d8	8.710	113	402561	16.155	ng/u1	-0.02
21) Nitrobenzene-d5	9.169	128	375637	32.581	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.892	143	355395	34.414	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.434	165	647815	27.068	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.951	131	590829	14.848	ng/u1	-0.01
46) Dimethylphthalate-d6	14.045	166	2206121	30.890	ng/u1	0.00
49) Acenaphthylene-d8	14.327	160	2590797	30.582	ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	66052	4.768	ng/u1	-0.02
60) Fluorene-d10	15.616	176	2048578	32.008	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.727	200	222634	27.953	ng/u1	-0.01
73) Anthracene-d10	17.463	188	3294194	35.790	ng/u1	-0.01
81) Pyrene-d10	19.751	212	3317646	49.440	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.809	264	2221833	37.735	ng/u1	-0.01

Target Compounds Qvalue

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

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