

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
 Data File : BM036727.D
 Acq On : 24 Sep 2022 23:36
 Operator : CG/JU
 Sample : N4662-05
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EXZ26

Quant Time: Sep 26 01:15:25 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	379741	20.000	ng/u1	0.00
20) Naphthalene-d8	10.816	136	1671561	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.627	164	1039091	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.368	188	1968360	20.000	ng/u1	0.00
79) Chrysene-d12	21.533	240	1316928	20.000	ng/u1	-0.01
88) Perylene-d12	23.962	264	1146361	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	38418	3.793	ng/uL	0.00
4) Pyridine-d5	3.816	84	226059	7.652	ng/u1	0.00
7) Phenol-d5	7.163	99	241421	7.114	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.334	67	624269	30.291	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	636162	26.038	ng/u1	0.00
15) 4-Methylphenol-d8	8.710	113	441980	17.626	ng/u1	-0.02
21) Nitrobenzene-d5	9.169	128	406202	35.903	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.892	143	383239	37.816	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.433	165	693758	29.539	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.951	131	722817	18.511	ng/u1	-0.01
46) Dimethylphthalate-d6	14.039	166	2295272	32.752	ng/u1	-0.01
49) Acenaphthylene-d8	14.327	160	2729381	32.834	ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	64689	4.759	ng/u1	-0.02
60) Fluorene-d10	15.615	176	2121878	33.786	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.727	200	216744	27.832	ng/u1	-0.01
73) Anthracene-d10	17.462	188	3320667	36.897	ng/u1	-0.01
81) Pyrene-d10	19.751	212	3284028	50.689	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.803	264	2196335	39.216	ng/u1	-0.02

Target Compounds Qvalue

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

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