

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
 Data File : BM036746.D
 Acq On : 26 Sep 2022 15:35
 Operator : CG/JU
 Sample : N4729-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EXFJ7

Quant Time: Sep 26 23:14:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 26 23:10:22 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.004	152	393107	20.000 ng/u1	0.00
20) Naphthalene-d8	10.816	136	1687918	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.627	164	1049798	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.368	188	2061449	20.000 ng/u1	0.00
79) Chrysene-d12	21.533	240	1551482	20.000 ng/u1	0.00
88) Perylene-d12	23.968	264	1331670	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.381	96	66216	6.315 ng/uL	0.00
4) Pyridine-d5	3.816	84	274766	8.984 ng/u1	0.00
7) Phenol-d5	7.157	99	436117	12.414 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	698630	32.747 ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	773270	30.574 ng/u1	0.00
15) 4-Methylphenol-d8	8.710	113	615822	23.724 ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	439461	38.466 ng/u1	0.00
24) 2-Nitrophenol-d4	9.892	143	413050	40.363 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.433	165	759294	32.017 ng/u1	0.00
31) 4-Chloroaniline-d4	10.951	131	946315	24.000 ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	2518147	35.565 ng/u1	0.00
49) Acenaphthylene-d8	14.327	160	2948974	35.113 ng/u1	0.00
54) 4-Nitrophenol-d4	14.815	143	126906	9.241 ng/u1	0.00
60) Fluorene-d10	15.615	176	2241696	35.330 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.733	200	242009	29.673 ng/u1	0.00
73) Anthracene-d10	17.468	188	3380089	35.861 ng/u1	0.00
81) Pyrene-d10	19.750	212	3583564	46.950 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.809	264	2427480	37.312 ng/u1	0.00

Target Compounds Qvalue

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

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