

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
 Data File : BM036592.D
 Acq On : 19 Sep 2022 22:24
 Operator : CG/JU
 Sample : N4604-07
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A5766

Quant Time: Sep 20 00:03:19 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 16 02:39:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	447040	20.000	ng/u1	0.00
20) Naphthalene-d8	10.839	136	1966322	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.651	164	1294905	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.386	188	2644411	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	2543357	20.000	ng/u1	-0.01
88) Perylene-d12	23.992	264	2348790	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	38277	3.210	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.175	99	728369	18.232	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	486985	20.073	ng/u1	0.00
11) 2-Chlorophenol-d4	7.551	132	598810	20.820	ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	574893	19.475	ng/u1	0.00
21) Nitrobenzene-d5	9.192	128	308030	23.144	ng/u1	0.00
24) 2-Nitrophenol-d4	9.916	143	286886	24.065	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	540543	19.565	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	349394	7.606	ng/u1	0.00
46) Dimethylphthalate-d6	14.063	166	1847213	21.151	ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	2238722	21.611	ng/u1	0.00
54) 4-Nitrophenol-d4	14.822	143	137509	8.118	ng/u1	0.00
60) Fluorene-d10	15.639	176	1693013	21.632	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	122434	11.702	ng/u1	0.00
73) Anthracene-d10	17.486	188	2623236	21.696	ng/u1	0.00
81) Pyrene-d10	19.762	212	2974656	23.774	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.833	264	2524442	21.999	ng/u1	0.00

Target Compounds Qvalue

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

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