

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
 Data File : BM036626.D
 Acq On : 21 Sep 2022 15:47
 Operator : CG/JU
 Sample : N4595-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Sep 21 22:50:00 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 21 22:42:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.022	152	348605	20.000	ng/u1	0.00
20) Naphthalene-d8	10.839	136	1560967	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.651	164	1023010	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.386	188	2106019	20.000	ng/u1	0.00
79) Chrysene-d12	21.550	240	1793218	20.000	ng/u1	0.00
88) Perylene-d12	23.997	264	1528812	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	39802	4.280	ng/uL	0.00
4) Pyridine-d5	3.828	84	187888	6.928	ng/u1	0.01
7) Phenol-d5	7.181	99	217793	6.991	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	569244	30.088	ng/u1	0.00
11) 2-Chlorophenol-d4	7.551	132	592685	26.425	ng/u1	0.00
15) 4-Methylphenol-d8	8.734	113	415461	18.049	ng/u1	0.00
21) Nitrobenzene-d5	9.192	128	394084	37.299	ng/u1	0.00
24) 2-Nitrophenol-d4	9.916	143	398285	42.086	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	702037	32.010	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	721619	19.790	ng/u1	0.00
46) Dimethylphthalate-d6	14.063	166	2695330	39.065	ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	3067705	37.484	ng/u1	0.00
54) 4-Nitrophenol-d4	14.833	143	96116	7.182	ng/u1	0.00
60) Fluorene-d10	15.639	176	2359991	38.168	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	344661	41.364	ng/u1	0.00
73) Anthracene-d10	17.486	188	3734705	38.785	ng/u1	0.00
81) Pyrene-d10	19.768	212	4182379	47.409	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.838	264	3076877	41.195	ng/u1	0.00

Target Compounds	Qvalue

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

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