

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
 Data File : BM036623.D
 Acq On : 21 Sep 2022 13:59
 Operator : CG/JU
 Sample : N4595-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Sep 21 22:49:00 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	345905	20.000	ng/u1	0.00
20) Naphthalene-d8	10.833	136	1591654	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.645	164	1069008	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.380	188	2174120	20.000	ng/u1	0.00
79) Chrysene-d12	21.550	240	1801343	20.000	ng/u1	0.00
88) Perylene-d12	23.991	264	1516039	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	43632	4.729	ng/uL	0.00
4) Pyridine-d5	3.822	84	172153	6.397	ng/u1	0.00
7) Phenol-d5	7.175	99	203989	6.599	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.345	67	572628	30.504	ng/u1	0.00
11) 2-Chlorophenol-d4	7.551	132	576427	25.901	ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	391046	17.121	ng/u1	0.00
21) Nitrobenzene-d5	9.186	128	390889	36.284	ng/u1	0.00
24) 2-Nitrophenol-d4	9.910	143	377067	39.075	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	667427	29.845	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	753332	20.261	ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	2560817	35.518	ng/u1	0.00
49) Acenaphthylene-d8	14.339	160	2927804	34.235	ng/u1	0.00
54) 4-Nitrophenol-d4	14.821	143	81372	5.819	ng/u1	0.00
60) Fluorene-d10	15.633	176	2272600	35.173	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	302001	35.109	ng/u1	0.00
73) Anthracene-d10	17.480	188	3630109	36.518	ng/u1	0.00
81) Pyrene-d10	19.762	212	3913179	44.157	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.833	264	2818081	38.048	ng/u1	0.00

Target Compounds Qvalue

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

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