

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112822\
 Data File : BM037766.D
 Acq On : 28 Nov 2022 19:06
 Operator : CG/JU
 Sample : N5710-12
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0KZ3

Quant Time: Nov 28 22:29:49 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 28 22:24:47 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.104	152	201114	20.000	ng/ul	0.00
20) Naphthalene-d8	10.928	136	929158	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.733	164	578366	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.468	188	1183924	20.000	ng/ul	0.00
79) Chrysene-d12	21.633	240	1115607	20.000	ng/ul	0.00
88) Perylene-d12	24.127	264	1001739	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	19160	3.951	ng/uL	0.00
4) Pyridine-d5	3.881	84	65112	4.125	ng/ul	0.01
7) Phenol-d5	7.257	99	408258	20.348	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.428	67	271702	20.300	ng/ul	0.00
11) 2-Chlorophenol-d4	7.634	132	305605	21.920	ng/ul	0.00
15) 4-Methylphenol-d8	8.810	113	322441	20.195	ng/ul	0.00
21) Nitrobenzene-d5	9.275	128	161495	22.171	ng/ul	0.00
24) 2-Nitrophenol-d4	10.004	143	162429	22.630	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.539	165	282610	21.516	ng/ul	0.00
31) 4-Chloroaniline-d4	11.063	131	285958	13.120	ng/ul	0.00
46) Dimethylphthalate-d6	14.139	166	884581	21.593	ng/ul	0.00
49) Acenaphthylene-d8	14.427	160	1056741	21.854	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	148595	16.467	ng/ul	0.00
60) Fluorene-d10	15.716	176	785525	21.894	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.827	200	90211	14.164	ng/ul	0.00
73) Anthracene-d10	17.568	188	1216219	21.968	ng/ul	0.00
81) Pyrene-d10	19.851	212	1339003	23.619	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.962	264	1129525	21.906	ng/ul	0.00

Target Compounds	Qvalue

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

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