

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
 Data File : BM037710.D
 Acq On : 26 Nov 2022 00:10
 Operator : CG/JU
 Sample : N5591-16
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C9828

Quant Time: Nov 26 04:14:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 24 01:27:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.110	152	151689	20.000	ng/u1	0.00
20) Naphthalene-d8	10.934	136	722969	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.739	164	448516	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.474	188	931650	20.000	ng/u1	0.00
79) Chrysene-d12	21.633	240	896429	20.000	ng/u1	0.00
88) Perylene-d12	24.127	264	884413	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	24612	6.729	ng/uL	0.00
4) Pyridine-d5	3.881	84	135969	11.421	ng/u1	0.01
7) Phenol-d5	7.257	99	482893	31.909	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.434	67	343013	33.978	ng/u1	0.00
11) 2-Chlorophenol-d4	7.634	132	354130	33.677	ng/u1	0.00
15) 4-Methylphenol-d8	8.816	113	381744	31.700	ng/u1	0.00
21) Nitrobenzene-d5	9.281	128	188608	33.278	ng/u1	0.00
24) 2-Nitrophenol-d4	10.010	143	178567	31.974	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.545	165	312539	30.581	ng/u1	0.00
31) 4-Chloroaniline-d4	11.069	131	514581	30.343	ng/u1	0.00
46) Dimethylphthalate-d6	14.139	166	1150641	36.219	ng/u1	0.00
49) Acenaphthylene-d8	14.433	160	1282393	34.198	ng/u1	0.00
54) 4-Nitrophenol-d4	14.916	143	209419	29.926	ng/u1	0.00
60) Fluorene-d10	15.721	176	987230	35.482	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.833	200	122214	24.386	ng/u1	0.00
73) Anthracene-d10	17.574	188	1594988	36.610	ng/u1	0.00
81) Pyrene-d10	19.851	212	1815495	39.853	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.968	264	1725673	37.908	ng/u1	0.00

Target Compounds Qvalue

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

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