

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
 Data File : BM037740.D
 Acq On : 26 Nov 2022 19:20
 Operator : CG/JU
 Sample : N5602-17
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AG9

Quant Time: Nov 27 21:29:36 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	231548	20.000	ng/u1	0.00
20) Naphthalene-d8	10.933	136	1076152	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.733	164	722895	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.474	188	1553636	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.639	240	1424856	20.000	ng/u1	# 0.00
88) Perylene-d12	24.127	264	1216991	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	23646	4.235	ng/uL	0.00
4) Pyridine-d5	3.881	84	134004	7.374	ng/u1	0.01
7) Phenol-d5	7.257	99	208923	9.044	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.428	67	541779	35.158	ng/u1	0.00
11) 2-Chlorophenol-d4	7.634	132	503969	31.397	ng/u1	0.00
15) 4-Methylphenol-d8	8.816	113	381992	20.780	ng/u1	0.00
21) Nitrobenzene-d5	9.281	128	360710	42.756	ng/u1	0.00
24) 2-Nitrophenol-d4	10.004	143	354552	42.651	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.545	165	609735	40.081	ng/u1	0.00
31) 4-Chloroaniline-d4	11.063	131	281831	11.164	ng/u1	-0.01
46) Dimethylphthalate-d6	14.145	166	2245567	43.855	ng/u1	0.00
49) Acenaphthylene-d8	14.433	160	2582990	42.738	ng/u1	0.00
54) 4-Nitrophenol-d4	14.921	143	100447	8.906	ng/u1	0.00
60) Fluorene-d10	15.721	176	2033393	45.343	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.833	200	329342	39.406	ng/u1	0.00
73) Anthracene-d10	17.574	188	3172537	43.667	ng/u1	0.00
81) Pyrene-d10	19.850	212	3726823	51.470	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.974	264	2897273	46.251	ng/u1	0.00

Target Compounds	Qvalue

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

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