

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022322\
 Data File : BM034051.D
 Acq On : 23 Feb 2022 14:11
 Operator : CG/JU
 Sample : N1634-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 20071

Quant Time: Feb 23 14:53:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 22 17:55:56 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.019	152	295865	20.000	ng	0.00
21) Naphthalene-d8	10.836	136	1008796	20.000	ng	# 0.00
39) Acenaphthene-d10	14.683	164	259387	20.000	ng	# 0.03
64) Phenanthrene-d10	17.530	188	170063	20.000	ng	# 0.14
76) Chrysene-d12	21.630	240	1360350	20.000	ng	# 0.06
86) Perylene-d12	24.047	264	1424910	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.566	112	1121873	67.594	ng	0.00
7) Phenol-d6	7.166	99	1454971	66.759	ng	0.00
23) Nitrobenzene-d5	9.178	82	1059029	56.653	ng	0.00
42) 2,4,6-Tribromophenol	16.207	330	311799	99.043	ng	0.07
45) 2-Fluorobiphenyl	13.289	172	2518256	129.346	ng	0.00
79) Terphenyl-d14	20.118	244	1249848	15.999	ng	0.11
Target Compounds						
						Qvalue
37) 2-Methylnaphthalene	12.495	142	80671	2.133	ng	# 27
46) 1,1'-Biphenyl	13.501	154	1283366	64.757	ng	84
50) Dimethylphthalate	14.030	163	452653	23.538	ng	# 55
66) n-Nitrosodiphenylamine	16.054	169	512064	96.177	ng	# 57
71) Phenanthrene	17.577	178	156574	16.598	ng	# 1
72) Anthracene	17.577	178	156574	17.056	ng	# 1

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

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