

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022322\
 Data File : BM034052.D
 Acq On : 23 Feb 2022 14:48
 Operator : CG/JU
 Sample : N1633-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 20220104-MR-7-(DISCRETE)

Quant Time: Feb 23 15:27:20 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	248073	20.000	ng	0.00
21) Naphthalene-d8	10.837	136	921556	20.000	ng	# 0.00
39) Acenaphthene-d10	14.654	164	556882	20.000	ng	0.00
64) Phenanthrene-d10	17.389	188	1107493	20.000	ng	# 0.00
76) Chrysene-d12	21.559	240	1121370	20.000	ng	0.00
86) Perylene-d12	24.024	264	1200824	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.572	112	2266656	162.879	ng	0.00
7) Phenol-d6	7.172	99	2726688	149.211	ng	0.00
23) Nitrobenzene-d5	9.184	82	1770925	103.703	ng	0.00
42) 2,4,6-Tribromophenol	16.136	330	1137571	168.311	ng	0.00
45) 2-Fluorobiphenyl	13.283	172	4441950	106.270	ng	0.00
79) Terphenyl-d14	20.012	244	6551287	101.733	ng	0.00

Target Compounds	Qvalue

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

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