

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022322\
 Data File : BM034062.D
 Acq On : 23 Feb 2022 20:55
 Operator : CG/JU
 Sample : N1634-01 10X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 20071

Quant Time: Feb 24 04:51:47 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	259745	20.000	ng	0.00
21) Naphthalene-d8	10.830	136	927541	20.000	ng	# 0.00
39) Acenaphthene-d10	14.654	164	619928	20.000	ng	0.00
64) Phenanthrene-d10	17.406	188	1236888	20.000	ng	# 0.01
76) Chrysene-d12	21.571	240	1196292	20.000	ng	0.00
86) Perylene-d12	24.024	264	1253846	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.566	112	93403	6.410	ng	0.00
7) Phenol-d6	7.160	99	117538	6.143	ng	-0.01
23) Nitrobenzene-d5	9.172	82	79073	4.601	ng	-0.01
42) 2,4,6-Tribromophenol	16.142	330	57341	7.621	ng	0.00
45) 2-Fluorobiphenyl	13.283	172	222534	4.783	ng	0.00
79) Terphenyl-d14	20.018	244	388351	5.653	ng	0.00
Target Compounds						
46) 1,1'-Biphenyl	13.489	154	103055	2.176	ng	Qvalue 93

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

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