

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011422\
 Data File : BM033734.D
 Acq On : 14 Jan 2022 18:48
 Operator : CG/JU
 Sample : N1157-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EFF-WW

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 01/17/2022
 Supervised By :Jagrut Upadhyay 01/17/2022

Quant Time: Jan 17 00:26:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM010522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 13 07:49:01 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.966	152	172643	20.000	ng	0.00
21) Naphthalene-d8	10.813	136	705645	20.000	ng	0.02
39) Acenaphthene-d10	14.607	164	469475	20.000	ng	0.00
64) Phenanthrene-d10	17.342	188	886939	20.000	ng	0.00
76) Chrysene-d12	21.500	240	617263	20.000	ng	0.00
86) Perylene-d12	23.918	264	691540	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.537	112	206889	18.900	ng	0.02
7) Phenol-d6	7.366	99	33835m	2.293	ng	0.24
23) Nitrobenzene-d5	9.201	82	1150518	75.298	ng	0.06
42) 2,4,6-Tribromophenol	16.089	330	515571	105.143	ng	0.00
45) 2-Fluorobiphenyl	13.242	172	2764495	77.126	ng	0.00
79) Terphenyl-d14	19.954	244	3040490	88.663	ng	0.00
Target Compounds						
32) Benzoic acid	10.360	122	19032	2.646	ng	Qvalue 95

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

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