

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF013122\
 Data File : BF126754.D
 Acq On : 31 Jan 2022 16:40
 Operator : CG/JU
 Sample : N1319-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-1S

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 02/01/2022
 Supervised By :Jagrut Upadhyay 02/01/2022

Quant Time: Feb 01 02:41:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 27 03:09:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.957	152	150458	20.000	ng	0.00
21) Naphthalene-d8	8.239	136	607795	20.000	ng	0.00
39) Acenaphthene-d10	9.998	164	342078	20.000	ng	0.00
64) Phenanthrene-d10	11.492	188	587614	20.000	ng	0.00
76) Chrysene-d12	14.139	240	394819	20.000	ng	# 0.00
86) Perylene-d12	15.663	264	295041	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.598	112	1252847	109.574	ng	0.03
7) Phenol-d6	6.581	99	1541049	97.007	ng	0.00
23) Nitrobenzene-d5	7.522	82	1322826	85.188	ng	0.00
42) 2,4,6-Tribromophenol	10.792	330	434683	135.782	ng	0.00
45) 2-Fluorobiphenyl	9.316	172	1732499	77.565	ng	0.00
79) Terphenyl-d14	13.086	244	1977260	83.998	ng	0.00
Target Compounds						
32) Benzoic acid	7.922	122	19523m	3.468	ng	Qvalue

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

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