

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012522\
 Data File : BF126667.D
 Acq On : 25 Jan 2022 16:56
 Operator : CG/JU
 Sample : N1250-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-3

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 01/26/2022
 Supervised By :Jagrut Upadhyay 01/26/2022

Quant Time: Jan 26 00:21:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 19 07:04:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.963	152	117312	20.000	ng	0.00
21) Naphthalene-d8	8.245	136	458813	20.000	ng	0.00
39) Acenaphthene-d10	10.004	164	249766	20.000	ng	0.00
64) Phenanthrene-d10	11.498	188	393790	20.000	ng	0.00
76) Chrysene-d12	14.145	240	233497	20.000	ng	# 0.00
86) Perylene-d12	15.668	264	233850	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	605410	67.909	ng	0.00
7) Phenol-d6	6.569	99	589021	47.554	ng	-0.01
23) Nitrobenzene-d5	7.528	82	1133257	96.678	ng	0.00
42) 2,4,6-Tribromophenol	10.798	330	318776	136.379	ng	0.00
45) 2-Fluorobiphenyl	9.322	172	1428202	87.573	ng	0.00
79) Terphenyl-d14	13.086	244	1300082	93.388	ng	0.00
Target Compounds						
32) Benzoic acid	7.951	122	42482m	9.997	ng	Qvalue
38) 1-Methylnaphthalene	9.057	142	129850	8.691	ng	99

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

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