

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101320\
 Data File : BF121942.D
 Acq On : 13 Oct 2020 17:13
 Operator : JU/CG
 Sample : L4333-12
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OH-WC-C-C1

Manual Integrations
 APPROVED

mohammad
 10/14/2020 12:53:31 PM

Quant Time: Oct 13 18:05:56 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 12 16:22:33 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.745	152	113524	20.00	ng	0.00
21) Naphthalene-d8	8.028	136	433051	20.00	ng	0.00
39) Acenaphthene-d10	9.775	164	212458	20.00	ng	-0.01
64) Phenanthrene-d10	11.263	188	364910	20.00	ng	-0.01
76) Chrysene-d12	13.898	240	265915	20.00	ng	-0.02
86) Perylene-d12	15.339	264	264999	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	783466	101.86	ng	0.01
7) Phenol-d6	6.393	99	852191	86.07	ng	0.00
23) Nitrobenzene-d5	7.310	82	799961	94.74	ng	-0.01
42) 2,4,6-Tribromophenol	10.569	330	326905	124.39	ng	-0.01
45) 2-Fluorobiphenyl	9.104	172	1283932	88.92	ng	0.00
79) Terphenyl-d14	12.851	244	1195099	85.65	ng	-0.01
Target Compounds						
32) Benzoic acid	7.840	122	9751m	9.71	ng	Qvalue

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

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