

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012422\
 Data File : BF126647.D
 Acq On : 24 Jan 2022 18:58
 Operator : CG/JU
 Sample : N1243-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AMI-1

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/25/2022
 Supervised By : Jagrut Upadhyay 01/25/2022

Quant Time: Jan 25 03:14:23 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	106593	20.000	ng	0.00
21) Naphthalene-d8	8.245	136	407804	20.000	ng	0.00
39) Acenaphthene-d10	10.004	164	223810	20.000	ng	0.00
64) Phenanthrene-d10	11.498	188	371314	20.000	ng	0.00
76) Chrysene-d12	14.139	240	234254	20.000	ng	#-0.01
86) Perylene-d12	15.662	264	224107	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	772857	95.410	ng	0.00
7) Phenol-d6	6.575	99	1056812	93.901	ng	0.00
23) Nitrobenzene-d5	7.522	82	792420	76.057	ng	-0.01
42) 2,4,6-Tribromophenol	10.792	330	214608	102.462	ng	0.00
45) 2-Fluorobiphenyl	9.322	172	996311	68.176	ng	0.00
79) Terphenyl-d14	13.086	244	921087	65.950	ng	0.00
Target Compounds						
32) Benzoic acid	7.928	122	19798m	5.242	ng	Qvalue

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

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