

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080422\
 Data File : BF129582.D
 Acq On : 04 Aug 2022 22:04
 Operator : CG\JU
 Sample : N3930-10
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RVE-102

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/05/2022
 Supervised By : Jagrut Upadhyay 08/05/2022

Quant Time: Aug 05 02:10:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 02:05:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.857	152	159234	20.00	ng	0.00
21) Naphthalene-d8	8.140	136	578876	20.00	ng	0.00
39) Acenaphthene-d10	9.892	164	252254	20.00	ng	-0.01
64) Phenanthrene-d10	11.386	188	364931	20.00	ng	0.00
76) Chrysene-d12	14.033	240	311597	20.00	ng	0.00
86) Perylene-d12	15.515	264	273520	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	1065834	109.04	ng	0.00
7) Phenol-d6	6.504	99	1292921	105.23	ng	-0.01
23) Nitrobenzene-d5	7.422	82	797993	75.25	ng	-0.01
42) 2,4,6-Tribromophenol	10.692	330	245510	101.23	ng	0.00
45) 2-Fluorobiphenyl	9.216	172	1304491	80.00	ng	0.00
79) Terphenyl-d14	12.969	244	1095439	66.32	ng	-0.01
Target Compounds						
75) Fluoranthene	12.598	202	49268	2.44	ng	97
78) Pyrene	12.833	202	62461	2.40	ng	99

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

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