

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080222\
 Data File : BF129521.D
 Acq On : 02 Aug 2022 22:41
 Operator : CG\JU
 Sample : N3927-13
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RVE-86

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/03/2022
 Supervised By : Jagrut Upadhyay 08/03/2022

Quant Time: Aug 03 01:48:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 01:11:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	151543	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	578579	20.000	ng	0.00
39) Acenaphthene-d10	9.916	164	269562	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	370877	20.000	ng	0.00
76) Chrysene-d12	14.045	240	278028	20.000	ng	0.00
86) Perylene-d12	15.533	264	218303	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	940587	101.108	ng	0.00
7) Phenol-d6	6.522	99	1177141	100.670	ng	0.00
23) Nitrobenzene-d5	7.440	82	751349	70.889	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	231253	89.230	ng	0.00
45) 2-Fluorobiphenyl	9.234	172	1295922	74.370	ng	0.00
79) Terphenyl-d14	12.986	244	989354	67.133	ng	0.00
Target Compounds						
78) Pyrene	12.845	202	65250	2.807	ng	Qvalue 98

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

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