

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080922\
 Data File : BF129684.D
 Acq On : 10 Aug 2022 01:05
 Operator : CG\JU
 Sample : N3962-14
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RVE-199D

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 08/10/2022
 Supervised By :Jagrut Upadhyay 08/10/2022

Quant Time: Aug 10 05:38:01 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 16:27:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.840	152	164353	20.000	ng	-0.01
21) Naphthalene-d8	8.122	136	619655	20.000	ng	-0.01
39) Acenaphthene-d10	9.880	164	285846	20.000	ng	-0.01
64) Phenanthrene-d10	11.369	188	413426	20.000	ng	-0.01
76) Chrysene-d12	14.015	240	299242	20.000	ng	-0.01
86) Perylene-d12	15.492	264	242593	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	925631	91.745	ng	0.00
7) Phenol-d6	6.492	99	1169769	92.242	ng	-0.01
23) Nitrobenzene-d5	7.404	82	726402	63.992	ng	-0.02
42) 2,4,6-Tribromophenol	10.675	330	240946	87.674	ng	-0.01
45) 2-Fluorobiphenyl	9.198	172	1272357	68.858	ng	-0.01
79) Terphenyl-d14	12.957	244	1055052	66.516	ng	-0.01
Target Compounds						
75) Fluoranthene	12.586	202	97777	4.276	ng	99
78) Pyrene	12.816	202	117330	4.689	ng	97
83) Chrysene	14.039	228	60792	3.156	ng	98
88) Benzo(b)fluoranthene	15.063	252	74891m	4.654	ng	
90) Benzo(a)pyrene	15.427	252	48015	3.676	ng	# 93

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

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