

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080522\
 Data File : BF129612.D
 Acq On : 05 Aug 2022 21:50
 Operator : CG\JU
 Sample : N3961-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RVE-169

Manual Integrations
 APPROVED

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

Quant Time: Aug 08 02:02:41 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	6.845	152	174814	20.00	ng	0.00
21) Naphthalene-d8	8.128	136	680924	20.00	ng	0.00
39) Acenaphthene-d10	9.880	164	325838	20.00	ng	-0.01
64) Phenanthrene-d10	11.369	188	461050	20.00	ng	#-0.01
76) Chrysene-d12	14.015	240	298955	20.00	ng	-0.01
86) Perylene-d12	15.492	264	330239	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	990529	92.30	ng	0.00
7) Phenol-d6	6.492	99	1240826	91.99	ng	-0.01
23) Nitrobenzene-d5	7.410	82	783103	62.78	ng	-0.01
42) 2,4,6-Tribromophenol	10.675	330	265798	84.85	ng	-0.01
45) 2-Fluorobiphenyl	9.204	172	1372256	65.15	ng	0.00
79) Terphenyl-d14	12.957	244	982907	62.03	ng	-0.01
Target Compounds						Qvalue
75) Fluoranthene	12.586	202	98086	3.85	ng	99
78) Pyrene	12.816	202	101456	4.06	ng	99
83) Chrysene	14.039	228	57385	2.98	ng	97
88) Benzo(b)fluoranthene	15.057	252	81176m	3.71	ng	
90) Benzo(a)pyrene	15.427	252	50963	2.87	ng	# 97

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

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