

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080922\
 Data File : BF129675.D
 Acq On : 09 Aug 2022 20:02
 Operator : CG\JU
 Sample : N3963-15
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RVE-219D

Manual Integrations
 APPROVED

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

Quant Time: Aug 10 05:33:56 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	178464	20.000	ng	-0.01
21) Naphthalene-d8	8.122	136	691006	20.000	ng	-0.01
39) Acenaphthene-d10	9.881	164	337944	20.000	ng	-0.01
64) Phenanthrene-d10	11.369	188	484285	20.000	ng	-0.01
76) Chrysene-d12	14.016	240	305403	20.000	ng	-0.01
86) Perylene-d12	15.492	264	312542	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	1134715	103.575	ng	0.00
7) Phenol-d6	6.493	99	1437571	104.396	ng	-0.01
23) Nitrobenzene-d5	7.410	82	901450	71.213	ng	-0.01
42) 2,4,6-Tribromophenol	10.675	330	325370	100.142	ng	-0.01
45) 2-Fluorobiphenyl	9.198	172	1591626	72.857	ng	-0.01
79) Terphenyl-d14	12.957	244	1221856	75.478	ng	-0.01
Target Compounds						Qvalue
75) Fluoranthene	12.580	202	115553	4.314	ng	98
78) Pyrene	12.816	202	110311	4.319	ng	95
81) Benzo(a)anthracene	14.004	228	57003	2.732	ng	96
83) Chrysene	14.039	228	56686	2.883	ng	97
88) Benzo(b)fluoranthene	15.057	252	95683m	4.615	ng	
90) Benzo(a)pyrene	15.427	252	60833	3.615	ng	# 96

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

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