

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061722\
 Data File : BF128860.D
 Acq On : 17 Jun 2022 18:10
 Operator : CG\JU
 Sample : N3367-01RE
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MJS-LB-02-13.0-14.0RE

Quant Time: Jun 18 05:07:12 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 14 14:53:53 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 06/20/2022
 Supervised By :Jagrut Upadhyay 06/20/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	53389	20.000	ng	0.00
21) Naphthalene-d8	8.063	136	207449	20.000	ng	# 0.00
39) Acenaphthene-d10	9.816	164	111329	20.000	ng	-0.01
64) Phenanthrene-d10	11.304	188	169915	20.000	ng	-0.01
76) Chrysene-d12	13.951	240	124259	20.000	ng	#-0.01
86) Perylene-d12	15.404	264	113719	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	527112	157.261	ng	0.00
7) Phenol-d6	6.434	99	682320	166.766	ng	0.00
23) Nitrobenzene-d5	7.345	82	479032	107.511	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	174457	132.626	ng	0.00
45) 2-Fluorobiphenyl	9.139	172	881389	110.901	ng	0.00
79) Terphenyl-d14	12.892	244	645988	90.335	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	8.081	128	21364	2.007	ng	98
71) Phenanthrene	11.327	178	53153	6.237	ng	98
75) Fluoranthene	12.522	202	45103	5.063	ng	95
78) Pyrene	12.751	202	38253	4.138	ng	100
81) Benzo(a)anthracene	13.939	228	19238	2.484	ng	98
83) Chrysene	13.974	228	16765	2.270	ng	95
88) Benzo(b)fluoranthene	14.986	252	19242m	2.648	ng	
90) Benzo(a)pyrene	15.345	252	14417	2.524	ng	# 96

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

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