

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061722\
 Data File : BF128861.D
 Acq On : 17 Jun 2022 18:40
 Operator : CG\JU
 Sample : N3364-01RE
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-02-061522RE

Manual Integrations
 APPROVED

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

Quant Time: Jun 18 05:07:19 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	63152	20.000	ng	0.00
21) Naphthalene-d8	8.063	136	247314	20.000	ng	0.00
39) Acenaphthene-d10	9.822	164	135700	20.000	ng	0.00
64) Phenanthrene-d10	11.304	188	176017	20.000	ng	-0.01
76) Chrysene-d12	13.957	240	149398	20.000	ng	# 0.00
86) Perylene-d12	15.415	264	106995	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	494350	124.686	ng	0.00
7) Phenol-d6	6.434	99	562670	116.262	ng	0.00
23) Nitrobenzene-d5	7.345	82	369177	69.500	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	133638	83.349	ng	0.00
45) 2-Fluorobiphenyl	9.139	172	695850	71.831	ng	0.00
79) Terphenyl-d14	12.898	244	597563	69.502	ng	0.00
Target Compounds						Qvalue
75) Fluoranthene	12.521	202	44256	4.796	ng	97
78) Pyrene	12.751	202	34822	3.133	ng	99
81) Benzo(a)anthracene	13.945	228	21624	2.322	ng	# 87
83) Chrysene	13.980	228	30809	3.470	ng	# 88
84) Bis(2-ethylhexyl)phtha...	13.933	149	15453	2.546	ng	# 95
87) Indeno(1,2,3-cd)pyrene	16.856	276	13196	2.046	ng	94
88) Benzo(b)fluoranthene	14.992	252	32472m	4.750	ng	
90) Benzo(a)pyrene	15.351	252	15389	2.864	ng	# 75
92) Benzo(g,h,i)perylene	17.303	276	14858	2.803	ng	# 85

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

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