

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061622\
 Data File : BF128846.D
 Acq On : 16 Jun 2022 18:31
 Operator : CG\JU
 Sample : N3353-01DL 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-01-061422DL

Quant Time: Jun 17 02:24:07 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	98293	20.000	ng	0.00
21) Naphthalene-d8	8.069	136	345450	20.000	ng	# 0.00
39) Acenaphthene-d10	9.822	164	147403	20.000	ng	0.00
64) Phenanthrene-d10	11.310	188	257577	20.000	ng	0.00
76) Chrysene-d12	13.963	240	150089	20.000	ng	# 0.00
86) Perylene-d12	15.421	264	114024	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	230279	37.317	ng	0.00
7) Phenol-d6	6.440	99	282823	37.546	ng	0.00
23) Nitrobenzene-d5	7.345	82	185759	25.036	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	54350	31.206	ng	0.00
45) 2-Fluorobiphenyl	9.139	172	350368	33.296	ng	0.00
79) Terphenyl-d14	12.898	244	268759	31.115	ng	0.00
Target Compounds						
						Qvalue
49) Acenaphthylene	9.686	152	33018	2.556	ng	96
71) Phenanthrene	11.333	178	132646	10.267	ng	99
72) Anthracene	11.386	178	48942	3.827	ng	97
75) Fluoranthene	12.533	202	439878	32.572	ng	94
78) Pyrene	12.757	202	414203	37.093	ng	98
81) Benzo(a)anthracene	13.951	228	185408	19.818	ng	96
83) Chrysene	13.986	228	185622m	20.811	ng	
87) Indeno(1,2,3-cd)pyrene	16.868	276	84162m	12.246	ng	
88) Benzo(b)fluoranthene	14.998	252	218220m	29.951	ng	
89) Benzo(k)fluoranthene	15.021	252	63941m	9.329	ng	
90) Benzo(a)pyrene	15.357	252	147527m	25.762	ng	
91) Dibenzo(a,h)anthracene	16.874	278	21075	3.697	ng	# 82
92) Benzo(g,h,i)perylene	17.309	276	78416	13.879	ng	# 90

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

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