

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041222\
 Data File : BF127765.D
 Acq On : 12 Apr 2022 18:31
 Operator : CG\JU
 Sample : N2364-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-03-COMP

Quant Time: Apr 13 07:25:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 07:19:18 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/13/2022
 Supervised By : Jagrut Upadhyay 04/13/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.957	152	116574	20.000	ng	0.00
21) Naphthalene-d8	8.239	136	425185	20.000	ng	# 0.00
39) Acenaphthene-d10	9.998	164	244217	20.000	ng	0.00
64) Phenanthrene-d10	11.486	188	447172	20.000	ng	0.00
76) Chrysene-d12	14.133	240	334547	20.000	ng	0.00
86) Perylene-d12	15.651	264	320952	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.575	112	705789	98.379	ng	0.00
7) Phenol-d6	6.575	99	895871	95.454	ng	0.00
23) Nitrobenzene-d5	7.516	82	639062	67.007	ng	-0.01
42) 2,4,6-Tribromophenol	10.786	330	287031	110.446	ng	0.00
45) 2-Fluorobiphenyl	9.316	172	1141872	66.145	ng	0.00
79) Terphenyl-d14	13.074	244	1232520	60.811	ng	0.00
Target Compounds						Qvalue
58) Fluorene	10.545	166	33560	2.188	ng	97
71) Phenanthrene	11.516	178	555813	23.459	ng	99
72) Anthracene	11.563	178	137188	5.854	ng	98
75) Fluoranthene	12.704	202	614835	23.781	ng	99
78) Pyrene	12.933	202	497858	17.623	ng	100
81) Benzo(a)anthracene	14.121	228	219445	9.821	ng	96
83) Chrysene	14.157	228	178946	8.343	ng	97
87) Indeno(1,2,3-cd)pyrene	17.198	276	83212	4.150	ng	96
88) Benzo(b)fluoranthene	15.198	252	187714m	9.092	ng	
89) Benzo(k)fluoranthene	15.227	252	78030m	3.810	ng	
90) Benzo(a)pyrene	15.586	252	142529	8.493	ng	98
92) Benzo(g,h,i)perylene	17.668	276	70660	4.242	ng	98

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

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