

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030422\
 Data File : BF127211.D
 Acq On : 04 Mar 2022 20:07
 Operator : CG\JU
 Sample : N1726-06DL 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 332DL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/07/2022
 Supervised By : Jagrut Upadhyay 03/07/2022

Quant Time: Mar 05 02:22:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 05 02:16:58 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	100946	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	401910	20.000	ng	# 0.00
39) Acenaphthene-d10	9.916	164	198468	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	345770	20.000	ng	# 0.00
76) Chrysene-d12	14.057	240	205456	20.000	ng	# 0.00
86) Perylene-d12	15.545	264	194191	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	308173	47.184	ng	0.00
7) Phenol-d6	6.504	99	399620	45.344	ng	-0.01
23) Nitrobenzene-d5	7.439	82	266254	29.902	ng	-0.01
42) 2,4,6-Tribromophenol	10.710	330	113647	49.734	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	446290	33.104	ng	0.00
79) Terphenyl-d14	12.998	244	448850	32.115	ng	0.00
Target Compounds						Qvalue
58) Fluorene	10.469	166	37398	2.834	ng	97
71) Phenanthrene	11.433	178	231693	11.972	ng	99
74) Di-n-butylphthalate	11.968	149	960007	42.265	ng	99
75) Fluoranthene	12.627	202	232579	12.657	ng	94
78) Pyrene	12.857	202	194494	10.896	ng	100
81) Benzo(a)anthracene	14.045	228	67387	4.865	ng	99
83) Chrysene	14.080	228	60154	4.618	ng	95
87) Indeno(1,2,3-cd)pyrene	17.045	276	29932	2.351	ng	# 84
88) Benzo(b)fluoranthene	15.104	252	66462m	5.096	ng	
89) Benzo(k)fluoranthene	15.133	252	26064m	2.072	ng	
90) Benzo(a)pyrene	15.480	252	47410m	4.654	ng	
92) Benzo(g,h,i)perylene	17.503	276	29670	2.858	ng	# 88

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

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