

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\
 Data File : BF119010.D
 Acq On : 21 Feb 2020 23:39
 Operator : CG/JU
 Sample : L1626-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-051-A

Manual Integrations
 APPROVED

mohammad
 2/24/2020 2:11:48 PM

Quant Time: Feb 24 03:43:16 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 20 17:40:17 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	142139	20.00	ng	-0.01
21) Naphthalene-d8	8.03	136	489792	20.00	ng	-0.01
39) Acenaphthene-d10	9.79	164	231136	20.00	ng	0.00
64) Phenanthrene-d10	11.27	188	403531	20.00	ng	0.00
76) Chrysene-d12	13.91	240	340401	20.00	ng	0.00
87) Perylene-d12	15.33	264	291710	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	813327	95.63	ng	0.00
7) Phenol-d6	6.40	99	923307	89.26	ng	0.00
23) Nitrobenzene-d5	7.32	82	596707	64.33	ng	-0.01
42) 2,4,6-Tribromophenol	10.57	330	257888	85.29	ng	0.00
45) 2-Fluorobiphenyl	9.11	172	1043094	66.48	ng	-0.01
79) Terphenyl-d14	12.85	244	1148501	58.09	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.50	163	105923	5.993	ng	99
71) Phenanthrene	11.29	178	131936	5.995	ng	98
75) Fluoranthene	12.48	202	196758	8.423	ng	99
78) Pyrene	12.71	202	210640	7.706	ng	100
81) Benzo(a)anthracene	13.90	228	107526	4.636	ng	95
83) Chrysene	13.93	228	103766	4.490	ng	96
88) Benzo(b)fluoranthene	14.93	252	115187m	5.991	ng	
90) Benzo(a)pyrene	15.27	252	74314	4.282	ng	# 96
92) Benzo(g,h,i)perylene	17.16	276	36950	2.449	ng	93

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

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