

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070519\
 Data File : BF115393.D
 Acq On : 5 Jul 2019 17:16
 Operator : HP/JU
 Sample : K3624-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-01-070319-B

Manual Integrations
 APPROVED

mohammad
 7/8/2019 2:12:58 PM

Quant Time: Jul 06 05:42:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 05 14:57:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	165766	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	655234	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	320482	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	555923	20.00	ng	0.00
76) Chrysene-d12	14.04	240	386587	20.00	ng	-0.02
87) Perylene-d12	15.52	264	416433	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	813488	75.82	ng	0.00
7) Phenol-d6	6.52	99	1071181	78.50	ng	0.00
23) Nitrobenzene-d5	7.45	82	655729	59.13	ng	0.00
42) 2,4,6-Tribromophenol	10.71	330	264324	74.59	ng	-0.01
45) 2-Fluorobiphenyl	9.25	172	1174817	64.33	ng	0.00
79) Terphenyl-d14	13.00	244	1095894	58.04	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.64	163	152983	6.360	ng	98
71) Phenanthrene	11.43	178	169033	5.781	ng	98
75) Fluoranthene	12.62	202	396714	12.887	ng	96
78) Pyrene	12.85	202	371714	12.365	ng	98
81) Benzo(a)anthracene	14.03	228	213589	8.623	ng	98
83) Chrysene	14.07	228	203841	8.009	ng	97
88) Benzo(b)fluoranthene	15.09	252	311568m	11.491	ng	
89) Benzo(k)fluoranthene	15.11	252	98897m	4.101	ng	
90) Benzo(a)pyrene	15.45	252	198755	8.499	ng	100
92) Benzo(g,h,i)perylene	17.43	276	107383	5.677	ng	98

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

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