

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110719\
 Data File : BF117718.D
 Acq On : 7 Nov 2019 21:21
 Operator : JU
 Sample : K5722-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 4-A

Manual Integrations
 APPROVED

Sohil
 11/8/2019 3:31:41 PM

Quant Time: Nov 08 02:44:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 17:27:10 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	318751	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	1219930	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	632463	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	1087299	20.00	ng	0.00
76) Chrysene-d12	14.12	240	722692	20.00	ng	0.00
87) Perylene-d12	15.63	264	895780	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	1838261	104.39	ng	0.00
7) Phenol-d6	6.55	99	2347221	108.10	ng	0.00
23) Nitrobenzene-d5	7.49	82	1412434	68.36	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	635947	102.79	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	2541341	79.28	ng	0.00
79) Terphenyl-d14	13.06	244	2390560	68.10	ng	0.00
Target Compounds						
10) Phenol	6.56	94	84257	3.601	ng	95
50) Dimethylphthalate	9.69	163	373792	8.742	ng	99
71) Phenanthrene	11.50	178	910987	17.480	ng	95
72) Anthracene	11.55	178	136975	2.642	ng	96
75) Fluoranthene	12.69	202	842341	15.829	ng	95
78) Pyrene	12.92	202	1125036	21.325	ng	96
81) Benzo(a)anthracene	14.11	228	486500	10.815	ng	98
83) Chrysene	14.14	228	517777	11.655	ng	98
86) Indeno(1,2,3-cd)pyrene	17.16	276	175801	4.252	ng	95
88) Benzo(b)fluoranthene	15.18	252	483625m	8.572	ng	
89) Benzo(k)fluoranthene	15.21	252	110378m	2.194	ng	
90) Benzo(a)pyrene	15.56	252	352248	7.202	ng	98
92) Benzo(a,h,i)perylene	17.62	276	176778	4.159	ng	99

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

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