

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110719\
 Data File : BF117720.D
 Acq On : 7 Nov 2019 22:15
 Operator : JU
 Sample : K5722-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 4-C

Manual Integrations
 APPROVED

Sohil
 11/8/2019 3:31:43 PM

Quant Time: Nov 08 02:51:01 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	309878	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	1190520	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	626785	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	1124761	20.00	ng	0.00
76) Chrysene-d12	14.12	240	763827	20.00	ng	0.00
87) Perylene-d12	15.63	264	901576	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	1539085	89.91	ng	0.00
7) Phenol-d6	6.55	99	1959653	92.84	ng	0.00
23) Nitrobenzene-d5	7.49	82	1166467	57.85	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	551309	89.92	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	2127853	66.98	ng	0.00
79) Terphenyl-d14	13.06	244	2161949	58.27	ng	0.00
Target Compounds						
10) Phenol	6.56	94	68947	3.031	ng	89
50) Dimethylphthalate	9.69	163	317893	7.502	ng	99
71) Phenanthrene	11.49	178	621902	11.536	ng	95
75) Fluoranthene	12.69	202	483176	8.777	ng	95
78) Pyrene	12.92	202	757558	13.586	ng	96
81) Benzo(a)anthracene	14.11	228	286331	6.022	ng	97
83) Chrysene	14.14	228	310743	6.618	ng	98
88) Benzo(b)fluoranthene	15.18	252	240103m	4.228	ng	
90) Benzo(a)pyrene	15.56	252	193925	3.940	ng	99
92) Benzo(g,h,i)perylene	17.62	276	93835	2.193	ng	98

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

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