

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110719\
 Data File : BF117721.D
 Acq On : 7 Nov 2019 22:42
 Operator : JU
 Sample : K5722-04
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 4-D

Manual Integrations
 APPROVED

Sohil
 11/8/2019 3:31:44 PM

Quant Time: Nov 08 02:53:03 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	313753	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	1210856	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	623800	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	1115001	20.00	ng	0.00
76) Chrysene-d12	14.12	240	732380	20.00	ng	0.00
87) Perylene-d12	15.63	264	891652	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	1830456	105.61	ng	0.00
7) Phenol-d6	6.55	99	2325619	108.82	ng	0.00
23) Nitrobenzene-d5	7.49	82	1406547	68.58	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	652870	106.99	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	2575050	81.45	ng	0.00
79) Terphenyl-d14	13.06	244	2540656	71.41	ng	0.00
Target Compounds						
10) Phenol	6.56	94	82408	3.578	ng	97
50) Dimethylphthalate	9.69	163	361869	8.581	ng	98
71) Phenanthrene	11.50	178	1012149	18.938	ng	95
72) Anthracene	11.54	178	142683	2.683	ng	95
75) Fluoranthene	12.69	202	810537	14.853	ng	95
78) Pyrene	12.92	202	1200960	22.463	ng	97
81) Benzo(a)anthracene	14.11	228	485736	10.655	ng	98
83) Chrysene	14.14	228	534246	11.867	ng	98
86) Indeno(1,2,3-cd)pyrene	17.16	276	173937	4.151	ng	96
88) Benzo(b)fluoranthene	15.18	252	442392m	7.878	ng	
89) Benzo(k)fluoranthene	15.21	252	143340m	2.862	ng	
90) Benzo(a)pyrene	15.56	252	354149	7.275	ng	99
92) Benzo(a,h,i)perylene	17.62	276	182625	4.316	ng	98

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

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