

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041819\
 Data File : BF113820.D
 Acq On : 18 Apr 2019 16:38
 Operator : JU/SJ
 Sample : K2411-11
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-14-D

Manual Integrations
 APPROVED

Sohil
 4/23/2019 8:16:41 AM

Quant Time: Apr 18 17:07:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 18 13:55:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	190309	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	696608	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	335006	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	579783	20.00	ng	0.00
76) Chrysene-d12	14.04	240	467698	20.00	ng	0.00
87) Perylene-d12	15.51	264	608311	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	1070470	93.53	ng	0.00
7) Phenol-d6	6.52	99	1355490	94.97	ng	0.00
23) Nitrobenzene-d5	7.45	82	824545	72.08	ng	0.00
42) 2,4,6-Tribromophenol	10.71	330	327537	88.49	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1518191	69.11	ng	0.00
79) Terphenyl-d14	12.99	244	1486777	66.66	ng	0.00
Target Compounds						
49) Acenaphthylene	9.79	152	80572	2.482	ng	97
50) Dimethylphthalate	9.63	163	160302	6.838	ng	100
71) Phenanthrene	11.43	178	91931	3.113	ng	100
75) Fluoranthene	12.62	202	115856	3.702	ng	98
78) Pyrene	12.85	202	176080	5.342	ng	97
81) Benzo(a)anthracene	14.03	228	171627m	6.278	ng	
83) Chrysene	14.06	228	191247m	7.048	ng	
86) Indeno(1,2,3-cd)pyrene	16.97	276	259863	11.731	ng	95
88) Benzo(b)fluoranthene	15.08	252	458971m	12.103	ng	
89) Benzo(k)fluoranthene	15.10	252	114845m	3.325	ng	
90) Benzo(a)pyrene	15.44	252	308537	9.201	ng	97
91) Dibenzo(a,h)anthracene	16.99	278	74846m	2.509	ng	
92) Benzo(a,h,i)perylene	17.42	276	235505	7.766	ng	99

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

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