

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111618\
 Data File : BF110817.D
 Acq On : 16 Nov 2018 15:31
 Operator : JU/SJ
 Sample : J5952-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VTW-01(7-10)

Manual Integrations
 APPROVED

Sohil
 11/19/2018 11:55:46 AM

Quant Time: Nov 16 16:27:57 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	66609	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	278327	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	130756	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	232577	20.00	ng	-0.01
76) Chrysene-d12	14.13	240	149088	20.00	ng	-0.01
87) Perylene-d12	15.66	264	146960	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	387206	94.42	ng	0.00
7) Phenol-d6	6.56	99	495822	94.55	ng	-0.01
23) Nitrobenzene-d5	7.50	82	307051	63.53	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	113854	86.41	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	546488	59.69	ng	-0.01
79) Terphenyl-d14	13.06	244	425882	53.50	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	8.25	128	80623	6.170	ng	98
37) 2-Methylnaphthalene	8.94	142	34103	3.982	ng	99
38) 1-Methylnaphthalene	9.03	142	30417m	3.693	ng	
49) Acenaphthylene	9.85	152	33071	2.613	ng	97
50) Dimethylphthalate	9.69	163	88465	9.067	ng	99
52) Acenaphthene	10.02	154	18594	2.325	ng	100
58) Fluorene	10.53	166	46168	5.137	ng	96
71) Phenanthrene	11.50	178	460760	36.978	ng	99
72) Anthracene	11.55	178	131476	10.460	ng	98
75) Fluoranthene	12.70	202	473997	37.943	ng	99
78) Pyrene	12.93	202	435387	37.928	ng	98
81) Benzo(a)anthracene	14.12	228	204026	22.896	ng	97
83) Chrysene	14.16	228	180732	21.288	ng	98
86) Indeno(1,2,3-cd)pyrene	17.23	276	77640	12.744	ng	98
88) Benzo(b)fluoranthene	15.20	252	215305m	24.490	ng	
89) Benzo(k)fluoranthene	15.23	252	75073m	8.642	ng	
90) Benzo(a)pyrene	15.59	252	166963	20.902	ng	97
91) Dibenzo(a,h)anthracene	17.23	278	21475m	3.177	ng	
92) Benzo(a,h,i)perylene	17.71	276	76048	11.339	ng	99

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

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