

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111518\
 Data File : BF110797.D
 Acq On : 15 Nov 2018 19:17
 Operator : JU/SJ
 Sample : J5767-11 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-111418-B

Manual Integrations
 APPROVED

Sohil
 11/16/2018 10:07:23 AM

Quant Time: Nov 16 07:21:55 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	63800	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	258223	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	110096	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	173550	20.00	ng	-0.01
76) Chrysene-d12	14.13	240	162299	20.00	ng	0.00
87) Perylene-d12	15.66	264	125056	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	199330	50.75	ng	-0.01
7) Phenol-d6	6.56	99	252605	50.29	ng	-0.01
23) Nitrobenzene-d5	7.50	82	156838	34.98	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	49185	44.34	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	276047	35.81	ng	-0.01
79) Terphenyl-d14	13.06	244	220927	25.49	ng	-0.01
Target Compounds						
10) Phenol	6.57	94	12737	2.187	ng	Qvalue 80
31) Naphthalene	8.25	128	224481	18.518	ng	99
37) 2-Methylnaphthalene	8.93	142	247257	31.115	ng	99
38) 1-Methylnaphthalene	9.03	142	213305m	27.911	ng	
46) 1,1'-Biphenyl	9.40	154	28632	2.788	ng	95
49) Acenaphthylene	9.85	152	39914	3.746	ng	# 90
50) Dimethylphthalate	9.69	163	27941	3.401	ng	100
52) Acenaphthene	10.02	154	28518	4.235	ng	92
58) Fluorene	10.53	166	115097	15.209	ng	# 98
71) Phenanthrene	11.50	178	452359	48.652	ng	99
72) Anthracene	11.56	178	84996	9.062	ng	96
75) Fluoranthene	12.70	202	232091	24.898	ng	98
78) Pyrene	12.93	202	396305	31.713	ng	99
81) Benzo(a)anthracene	14.12	228	162201m	16.720	ng	
83) Chrysene	14.16	228	186175m	20.145	ng	
86) Indeno(1,2,3-cd)pyrene	17.24	276	42796	6.453	ng	98
88) Benzo(b)fluoranthene	15.20	252	137360m	18.360	ng	
89) Benzo(k)fluoranthene	15.23	252	33552m	4.539	ng	
90) Benzo(a)pyrene	15.60	252	111648	16.425	ng	99
91) Dibenzo(a,h)anthracene	17.23	278	14016	2.437	ng	# 70
92) Benzo(g,h,i)perylene	17.72	276	42780	7.496	ng	96

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

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