

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111218\
 Data File : BF110692.D
 Acq On : 12 Nov 2018 23:22
 Operator : JU/SJ
 Sample : J5907-23 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-4-S

Manual Integrations
 APPROVED

Sohil
 11/13/2018 2:22:57 PM

Quant Time: Nov 13 06:43:08 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 09 15:13:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	106851	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	441605	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	193932	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	309144	20.00	ng	0.00
76) Chrysene-d12	14.13	240	197488	20.00	ng	-0.01
87) Perylene-d12	15.66	264	174750	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	362347	55.08	ng	0.00
7) Phenol-d6	6.57	99	474256	56.38	ng	0.00
23) Nitrobenzene-d5	7.51	82	299545	39.06	ng	0.00
42) 2,4,6-Tribromophenol	10.77	330	79070	40.46	ng	-0.01
45) 2-Fluorobiphenyl	9.30	172	536799	39.53	ng	0.00
79) Terphenyl-d14	13.06	244	349208	33.12	ng	-0.01
Target Compounds						
10) Phenol	6.58	94	26923	2.760	ng	Qvalue # 69
50) Dimethylphthalate	9.69	163	76597	5.293	ng	99
71) Phenanthrene	11.50	178	234709	14.171	ng	98
72) Anthracene	11.56	178	42186	2.525	ng	98
75) Fluoranthene	12.70	202	273410	16.466	ng	99
78) Pyrene	12.93	202	262217	17.244	ng	98
81) Benzo(a)anthracene	14.12	228	117965	9.994	ng	99
83) Chrysene	14.15	228	113564	10.098	ng	97
86) Indeno(1,2,3-cd)pyrene	17.22	276	45071	5.585	ng	96
88) Benzo(b)fluoranthene	15.20	252	135306m	12.943	ng	
89) Benzo(k)fluoranthene	15.23	252	35333m	3.420	ng	
90) Benzo(a)pyrene	15.59	252	86304	9.086	ng	97
92) Benzo(a,h,i)perylene	17.70	276	43686	5.478	ng	96

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

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