

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\
 Data File : BF110731.D
 Acq On : 13 Nov 2018 20:49
 Operator : JU/SJ
 Sample : J5908-07 2X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SR-3-S

Manual Integrations
 APPROVED

Sohil
 11/14/2018 1:52:25 PM

Quant Time: Nov 14 05:58:58 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	6.95	152	95185	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	385555	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	155096	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	226156	20.00	ng	0.00
76) Chrysene-d12	14.13	240	190127	20.00	ng	-0.01
87) Perylene-d12	15.66	264	147037	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	334022	57.00	ng	0.00
7) Phenol-d6	6.56	99	415059	55.39	ng	-0.01
23) Nitrobenzene-d5	7.50	82	258867	38.67	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	68506	43.84	ng	-0.01
45) 2-Fluorobiphenyl	9.30	172	448510	41.30	ng	0.00
79) Terphenyl-d14	13.06	244	291822	28.74	ng	-0.01
Target Compounds						
9) Benzaldehyde	6.56	77	680	Below Cal		Qvalue # 1
10) Phenol	6.58	94	19565	2.252 ng		# 72
50) Dimethylphthalate	9.69	163	47085	4.069 ng		98
71) Phenanthrene	11.50	178	36247	2.992 ng		97
75) Fluoranthene	12.70	202	117051	9.636 ng		98
78) Pyrene	12.93	202	105107	7.180 ng		98
81) Benzo(a)anthracene	14.12	228	71819	6.320 ng		99
83) Chrysene	14.16	228	64275	5.937 ng		97
86) Indeno(1,2,3-cd)pyrene	17.22	276	32109	4.133 ng		98
88) Benzo(b)fluoranthene	15.20	252	96278m	10.945 ng		
89) Benzo(k)fluoranthene	15.23	252	26161m	3.010 ng		
90) Benzo(a)pyrene	15.59	252	57403	7.183 ng		98
92) Benzo(a,h,i)perylene	17.70	276	28998	4.321 ng		99

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

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