

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121318\
 Data File : BF111391.D
 Acq On : 14 Dec 2018 1:05
 Operator : JU/SJ
 Sample : J6208-09
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3

Manual Integrations
 APPROVED

Sohil
 12/14/2018 5:11:29 PM

Quant Time: Dec 14 04:08:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 13:16:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	302854	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	1228636	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	509574	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	751916	20.00	ng	0.00
76) Chrysene-d12	13.95	240	521473	20.00	ng	0.00
87) Perylene-d12	15.39	264	412350	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	1413692	78.94	ng	0.02
7) Phenol-d6	6.45	99	1792212	74.22	ng	0.01
23) Nitrobenzene-d5	7.36	82	1271003	58.31	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	321817	71.64	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1782725	56.17	ng	0.00
79) Terphenyl-d14	12.90	244	1207152	50.95	ng	0.00
Target Compounds						
10) Phenol	6.45	94	113416	4.146	ng	95
50) Dimethylphthalate	9.55	163	151971	4.147	ng	97
71) Phenanthrene	11.33	178	229398	5.902	ng	94
75) Fluoranthene	12.52	202	381730	10.061	ng	98
78) Pyrene	12.75	202	355719	8.839	ng	94
81) Benzo(a)anthracene	13.94	228	159551	5.468	ng	97
83) Chrysene	13.97	228	159444	5.356	ng	95
86) Indeno(1,2,3-cd)pyrene	16.79	276	51682	2.220	ng	# 91
88) Benzo(b)fluoranthene	14.97	252	143977m	6.140	ng	
89) Benzo(k)fluoranthene	15.00	252	59002m	2.521	ng	
90) Benzo(a)pyrene	15.32	252	107422	4.920	ng	98
92) Benzo(a,h,i)perylene	17.22	276	49178	2.708	ng	98

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

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