

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091317\
 Data File : BF098507.D
 Acq On : 13 Sep 2017 23:32
 Operator : SJ/JU
 Sample : I5160-04
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G-7-1

Manual Integrations
 APPROVED

Sohil
 9/14/2017 5:52:57 PM

Quant Time: Sep 14 07:34:17 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF091117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 02:14:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	165392	20.00	ng	-0.01
21) Naphthalene-d8	7.94	136	666075	20.00	ng	-0.01
38) Acenaphthene-d10	9.69	164	274275	20.00	ng	-0.02
63) Phenanthrene-d10	11.17	188	409919	20.00	ng	-0.02
75) Chrysene-d12	13.80	240	340252	20.00	ng	-0.01
86) Perylene-d12	15.20	264	295622	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	1378881	123.27	ng	0.00
7) Phenol-d6	6.32	99	1650961	116.24	ng	0.00
23) Nitrobenzene-d5	7.23	82	929474	77.80	ng	-0.01
41) 2,4,6-Tribromophenol	10.48	330	211358	115.46	ng	-0.01
44) 2-Fluorobiphenyl	9.02	172	1388903	85.83	ng	-0.02
78) Terphenyl-d14	12.75	244	851789	54.04	ng	-0.02
Target Compounds						
10) Phenol	6.32	94	52557	3.186	ng	Qvalue # 66
49) Dimethylphthalate	9.41	163	259357	12.216	ng	98
70) Phenanthrene	11.19	178	145669	6.703	ng	98
71) Anthracene	11.25	178	58306	2.613	ng	97
74) Fluoranthene	12.38	202	215674	9.914	ng	98
77) Pyrene	12.60	202	213466	7.953	ng	98
80) Benzo(a)anthracene	13.79	228	119768m	6.004	ng	
82) Chrysene	13.83	228	121213m	6.011	ng	
85) Indeno(1,2,3-cd)pyrene	16.54	276	49123	3.472	ng	92
87) Benzo(b)fluoranthene	14.81	252	175169m	9.424	ng	
88) Benzo(k)fluoranthene	14.83	252	43005m	2.478	ng	
89) Benzo(a)pyrene	15.14	252	99670	6.019	ng	99
91) Benzo(a,h,i)perylene	16.94	276	43792	3.614	ng	98

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

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