

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091417\
 Data File : BF098545.D
 Acq On : 14 Sep 2017 23:18
 Operator : SJ/JU
 Sample : I5161-13
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RL-12-0-13

Manual Integrations
 APPROVED

Sohil
 9/15/2017 5:44:22 PM

Quant Time: Sep 15 06:49:18 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	6.65	152	133831	20.00	ng	-0.02
21) Naphthalene-d8	7.93	136	563629	20.00	ng	-0.02
38) Acenaphthene-d10	9.68	164	239285	20.00	ng	-0.02
63) Phenanthrene-d10	11.16	188	365028	20.00	ng	-0.02
75) Chrysene-d12	13.80	240	257302	20.00	ng	-0.02
86) Perylene-d12	15.19	264	290023	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	938307	103.66	ng	-0.02
7) Phenol-d6	6.31	99	1188520	103.42	ng	-0.01
23) Nitrobenzene-d5	7.22	82	701084	69.35	ng	-0.02
41) 2,4,6-Tribromophenol	10.47	330	150852	94.46	ng	-0.02
44) 2-Fluorobiphenyl	9.00	172	811583	57.49	ng	-0.03
78) Terphenyl-d14	12.74	244	531201	44.56	ng	-0.02
Target Compounds						
10) Phenol	6.32	94	37131	2.782	ng	# 61
20) 3+4-Methylphenols	7.10	107	27054	2.641	ng	# 57
49) Dimethylphthalate	9.40	163	191315	10.329	ng	99
70) Phenanthrene	11.19	178	188857	9.759	ng	97
74) Fluoranthene	12.37	202	321870	16.615	ng	97
77) Pyrene	12.60	202	267530	13.180	ng	98
80) Benzo(a)anthracene	13.79	228	140627	9.322	ng	98
82) Chrysene	13.82	228	145597	9.548	ng	97
85) Indeno(1,2,3-cd)pyrene	16.53	276	77324	7.228	ng	92
87) Benzo(b)fluoranthene	14.80	252	196758m	10.790	ng	
88) Benzo(k)fluoranthene	14.82	252	74063m	4.350	ng	
89) Benzo(a)pyrene	15.14	252	147327	9.068	ng	99
91) Benzo(a,h,i)perylene	16.93	276	70002	5.889	ng	93

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

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