

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091217\
 Data File : BF098467.D
 Acq On : 12 Sep 2017 23:06
 Operator : SJ/JU
 Sample : I5140-11
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RL-10-10.5-21

Manual Integrations
 APPROVED

Sohil
 9/13/2017 5:00:20 PM

Quant Time: Sep 13 02:23:39 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	206172	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	816745	20.00	ng	0.00
38) Acenaphthene-d10	9.69	164	311776	20.00	ng	-0.01
63) Phenanthrene-d10	11.17	188	477748	20.00	ng	-0.01
75) Chrysene-d12	13.81	240	397100	20.00	ng	0.00
86) Perylene-d12	15.21	264	307930	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	1506464	108.03	ng	0.00
7) Phenol-d6	6.32	99	1854090	104.72	ng	0.00
23) Nitrobenzene-d5	7.23	82	1063770	72.61	ng	0.00
41) 2,4,6-Tribromophenol	10.49	330	218911	105.20	ng	0.00
44) 2-Fluorobiphenyl	9.02	172	1503760	81.75	ng	-0.01
78) Terphenyl-d14	12.76	244	930563	50.58	ng	-0.01
Target Compounds						
10) Phenol	6.33	94	53774	2.615	ng	Qvalue 90
20) 3+4-Methylphenols	7.09	107	47388	3.003	ng	# 57
31) Naphthalene	7.97	128	80939	2.005	ng	98
49) Dimethylphthalate	9.42	163	231457	9.590	ng	99
74) Fluoranthene	12.39	202	71435	2.817	ng	98
77) Pyrene	12.61	202	69889	2.231	ng	97
80) Benzo(a)anthracene	13.80	228	47610m	2.045	ng	
82) Chrysene	13.83	228	47219	2.006	ng	96
83) Bis(2-ethylhexyl)phthalate	13.79	149	77087	4.520	ng	99
87) Benzo(b)fluoranthene	14.82	252	47960m	2.477	ng	
89) Benzo(a)pyrene	15.15	252	35123	2.036	ng	# 85

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

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