

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091417\
 Data File : BF098543.D
 Acq On : 14 Sep 2017 22:21
 Operator : SJ/JU
 Sample : I5161-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RL-14-14.25-28.5

Manual Integrations
 APPROVED

Sohil
 9/15/2017 5:44:00 PM

Quant Time: Sep 15 06:42:45 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	127606	20.00	ng	-0.02
21) Naphthalene-d8	7.93	136	533442	20.00	ng	-0.02
38) Acenaphthene-d10	9.68	164	235384	20.00	ng	-0.02
63) Phenanthrene-d10	11.16	188	379143	20.00	ng	-0.02
75) Chrysene-d12	13.79	240	267664	20.00	ng	-0.02
86) Perylene-d12	15.19	264	295885	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	996411	115.45	ng	-0.02
7) Phenol-d6	6.30	99	1223012	111.61	ng	-0.02
23) Nitrobenzene-d5	7.22	82	676447	70.69	ng	-0.02
41) 2,4,6-Tribromophenol	10.47	330	186811	118.91	ng	-0.02
44) 2-Fluorobiphenyl	9.01	172	917494	66.07	ng	-0.02
78) Terphenyl-d14	12.75	244	605128	48.80	ng	-0.02
Target Compounds						
10) Phenol	6.32	94	38668	3.038	ng	98
20) 3+4-Methylphenols	7.09	107	51253	5.248	ng	# 56
49) Dimethylphthalate	9.40	163	151689	8.325	ng	98
70) Phenanthrene	11.19	178	284717	14.165	ng	97
71) Anthracene	11.24	178	78101	3.785	ng	96
74) Fluoranthene	12.37	202	338778	16.837	ng	99
77) Pyrene	12.60	202	268171	12.700	ng	98
80) Benzo(a)anthracene	13.79	228	150123	9.566	ng	95
82) Chrysene	13.82	228	137495	8.668	ng	95
85) Indeno(1,2,3-cd)pyrene	16.53	276	61487	5.525	ng	93
87) Benzo(b)fluoranthene	14.80	252	163547m	8.791	ng	
88) Benzo(k)fluoranthene	14.83	252	60991m	3.512	ng	
89) Benzo(a)pyrene	15.13	252	127504	7.693	ng	96
91) Benzo(a,h,i)perylene	16.93	276	49703	4.099	ng	94

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

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