

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
 Data File : BF101694.D
 Acq On : 23 Dec 2017 10:50
 Operator : SJ/JU
 Sample : I6841-03
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-01-01

Manual Integrations
 APPROVED

Sohil
 12/26/2017 1:37:58 PM

Quant Time: Dec 25 02:45:25 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 22 16:07:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	148509	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	559794	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	254782	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	454801	20.00	ng	0.00
75) Chrysene-d12	13.96	240	306239	20.00	ng	0.00
86) Perylene-d12	15.43	264	272061	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1252265	125.61	ng	0.02
7) Phenol-d6	6.44	99	1423229	114.70	ng	0.00
23) Nitrobenzene-d5	7.36	82	983605	97.81	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	346414	141.10	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1598952	97.35	ng	0.00
78) Terphenyl-d14	12.90	244	1392650	109.63	ng	0.00
Target Compounds						
10) Phenol	6.45	94	415604	29.388	ng	99
17) 2-Methylphenol	7.05	107	126943	13.159	ng	98
20) 3+4-Methylphenols	7.21	107	367037	35.903	ng	88
27) 2,4-Dimethylphenol	7.73	122	104571	12.873	ng	97
31) Naphthalene	8.10	128	3965600	140.713	ng	98
37) 2-Methylnaphthalene	8.78	142	313002	17.047	ng	99
45) 1,1'-Biphenyl	9.25	154	70588	3.124	ng	97
48) Acenaphthylene	9.69	152	82753	3.030	ng	97
51) Acenaphthene	9.86	154	89450	5.452	ng	99
54) Dibenzofuran	10.03	168	155335	7.450	ng	98
57) Fluorene	10.37	166	164272	9.314	ng	97
70) Phenanthrene	11.34	178	235368	9.306	ng	99
72) Carbazole	11.55	167	526416	21.763	ng	99

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

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