

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
 Data File : BF101691.D
 Acq On : 23 Dec 2017 9:29
 Operator : SJ/JU
 Sample : I6838-03
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-7317-01

Manual Integrations
 APPROVED

Sohil
 12/26/2017 1:37:54 PM

Quant Time: Dec 26 10:58:17 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	157921	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	546344	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	266262	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	463022	20.00	ng	0.00
75) Chrysene-d12	13.96	240	309311	20.00	ng	0.00
86) Perylene-d12	15.43	264	260073	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.43	112	1326724	125.14	ng	0.02
7) Phenol-d6	6.44	99	1322218	100.21	ng	0.00
23) Nitrobenzene-d5	7.36	82	1027754	104.72	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	344386	134.23	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1573108	91.65	ng	0.00
78) Terphenyl-d14	12.90	244	1372122	106.94	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Pyridine	3.62	79	137619m	9.093	ng	
10) Phenol	6.49	94	3489712	232.054	ng	97
17) 2-Methylphenol	7.06	107	623963	60.824	ng	100
20) 3+4-Methylphenols	7.22	107	1905944	175.327	ng	88
27) 2,4-Dimethylphenol	7.74	122	418942	52.843	ng	98
31) Naphthalene	8.10	128	5947977	216.251	ng	98
37) 2-Methylnaphthalene	8.78	142	335576	18.726	ng	98
45) 1,1'-Biphenyl	9.25	154	78636	3.330	ng	97
48) Acenaphthylene	9.69	152	307512	10.776	ng	99
51) Acenaphthene	9.86	154	41446	2.417	ng	98
54) Dibenzofuran	10.03	168	208185	9.555	ng	99
57) Fluorene	10.37	166	166840	9.051	ng	96
70) Phenanthrene	11.34	178	236924	9.201	ng	99
72) Carbazole	11.55	167	488094	19.821	ng	99

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

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