

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122617\
 Data File : BF101713.D
 Acq On : 26 Dec 2017 14:12
 Operator : SJ/JU
 Sample : I6839-03DL 5X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-41446854-01DL

Manual Integrations
 APPROVED

Sohil
 12/27/2017 4:14:30 PM

Quant Time: Dec 26 15:17:22 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 26 10:48:01 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	146435	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	553674	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	232185	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	384271	20.00	ng	0.00
75) Chrysene-d12	13.97	240	300611	20.00	ng	0.00
86) Perylene-d12	15.43	264	281441	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	284395	28.93	ng	0.01
7) Phenol-d6	6.43	99	323110	26.41	ng	0.00
23) Nitrobenzene-d5	7.36	82	222440	22.36	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	67958	30.38	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	428192	28.61	ng	0.00
78) Terphenyl-d14	12.89	244	338450	27.14	ng	0.00
Target Compounds						
10) Phenol	6.45	94	523068	37.510	ng	99
17) 2-Methylphenol	7.05	107	111666	11.739	ng	98
20) 3+4-Methylphenols	7.21	107	360153	35.729	ng	87
27) 2,4-Dimethylphenol	7.75	122	65849	8.196	ng	96
31) Naphthalene	8.10	128	2126894	76.304	ng	99
37) 2-Methylnaphthalene	8.79	142	68163	3.753	ng	98
54) Dibenzofuran	10.03	168	41184	2.168	ng	99
70) Phenanthrene	11.34	178	43751	2.047	ng	98
72) Carbazole	11.56	167	138270	6.766	ng	99

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

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