

Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
 Data File : BF102158.D
 Acq On : 17 Jan 2018 19:00
 Operator : SJ/JU
 Sample : J1156-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-COMP

Manual Integrations
 APPROVED

Sohil
 1/18/2018 11:12:26 AM

Quant Time: Jan 18 03:04:50 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 17 15:28:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	156493	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	625501	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	267503	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	396293	20.00	ng	0.00
75) Chrysene-d12	13.86	240	267583	20.00	ng	-0.01
86) Perylene-d12	15.28	264	282402	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	1095223	98.83	ng	0.01
7) Phenol-d6	6.36	99	1267593	95.87	ng	0.00
23) Nitrobenzene-d5	7.29	82	807213	75.84	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	251474	107.72	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	1352478	78.99	ng	0.00
78) Terphenyl-d14	12.82	244	914114	70.93	ng	0.00
Target Compounds						
10) Phenol	6.37	94	58535	3.917	ng	89
49) Dimethylphthalate	9.47	163	153812	7.168	ng	99
74) Fluoranthene	12.44	202	91742	4.028	ng	95
77) Pyrene	12.67	202	81041	3.426	ng	99
80) Benzo(a)anthracene	13.86	228	39659	2.323	ng	99
82) Chrysene	13.89	228	39747	2.233	ng	97
85) Indeno(1,2,3-cd)pyrene	16.64	276	26232	2.025	ng	99
87) Benzo(b)fluoranthene	14.87	252	59127m	3.211	ng	
89) Benzo(a)pyrene	15.22	252	39006	2.354	ng	98

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

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