

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102377.D
 Acq On : 24 Jan 2018 8:23
 Operator : SJ/JU
 Sample : J1236-01 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CA-3

Manual Integrations
 APPROVED

Sohil
 1/24/2018 2:47:30 PM

Quant Time: Jan 24 10:38:05 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 23 00:43:41 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	126656	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	534531	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	231826	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	354679	20.00	ng	0.00
75) Chrysene-d12	13.87	240	255789	20.00	ng	0.00
86) Perylene-d12	15.29	264	232514	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	173836	20.81	ng	0.00
7) Phenol-d6	6.35	99	224897	22.33	ng	-0.01
23) Nitrobenzene-d5	7.27	82	131616	14.15	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	36430	15.86	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	255141	16.18	ng	-0.01
78) Terphenyl-d14	12.82	244	168791	14.88	ng	0.00
Target Compounds						
70) Phenanthrene	11.26	178	84942	4.110	ng	99
74) Fluoranthene	12.45	202	115914	5.500	ng	95
77) Pyrene	12.67	202	97295	4.920	ng	97
80) Benzo(a)anthracene	13.86	228	42692	2.711	ng	97
82) Chrysene	13.90	228	44208	2.764	ng	93
87) Benzo(b)fluoranthene	14.89	252	46756m	3.103	ng	
89) Benzo(a)pyrene	15.23	252	31777	2.338	ng	98

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

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