

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
 Data File : BF106441.D
 Acq On : 14 Jun 2018 14:49
 Operator : JU/SJ
 Sample : J3239-13
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-06-061318-C

Manual Integrations
 APPROVED

Sohil
 6/15/2018 11:16:46 AM

Quant Time: Jun 14 18:56:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 14 18:39:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	141142	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	536665	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	243318	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	411040	20.00	ng	0.00
75) Chrysene-d12	14.15	240	358935	20.00	ng	0.00
86) Perylene-d12	15.72	264	336414	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	755069	90.54	ng	0.00
7) Phenol-d6	6.61	99	944394	89.64	ng	0.00
23) Nitrobenzene-d5	7.52	82	640413	70.20	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	238178	101.74	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	1025773	70.90	ng	0.00
78) Terphenyl-d14	13.08	244	865536	59.89	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.71	163	149941	8.625	ng	99
70) Phenanthrene	11.52	178	60190	2.990	ng	97
74) Fluoranthene	12.71	202	102304	4.765	ng	96
77) Pyrene	12.94	202	93970	4.182	ng	98
80) Benzo(a)anthracene	14.14	228	59827	2.801	ng	97
82) Chrysene	14.18	228	53831	2.760	ng	96
87) Benzo(b)fluoranthene	15.26	252	67911m	3.339	ng	
89) Benzo(a)pyrene	15.65	252	45569	2.491	ng	94

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

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