

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
 Data File : BF106447.D
 Acq On : 14 Jun 2018 17:30
 Operator : JU/SJ
 Sample : J3241-07 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-05-061318-A

Manual Integrations
 APPROVED

Sohil
 6/15/2018 11:16:49 AM

Quant Time: Jun 14 19:06:59 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	117343	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	447148	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	183238	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	300578	20.00	ng	0.00
75) Chrysene-d12	14.15	240	344286	20.00	ng	0.00
86) Perylene-d12	15.70	264	266916	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	165289	23.84	ng	0.00
7) Phenol-d6	6.61	99	212273	24.23	ng	0.00
23) Nitrobenzene-d5	7.52	82	124400	16.37	ng	-0.02
41) 2,4,6-Tribromophenol	10.80	330	35356	20.05	ng	-0.01
44) 2-Fluorobiphenyl	9.32	172	232506	12.64	ng	0.00
78) Terphenyl-d14	13.08	244	218788	15.78	ng	-0.01
Target Compounds						
70) Phenanthrene	11.52	178	38630	2.624	ng	97
74) Fluoranthene	12.71	202	70624	4.499	ng	97
77) Pyrene	12.94	202	69993	3.247	ng	98
80) Benzo(a)anthracene	14.14	228	42417	2.070	ng	97
87) Benzo(b)fluoranthene	15.24	252	44159m	2.737	ng	

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

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