

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061318\
 Data File : BF106397.D
 Acq On : 13 Jun 2018 17:59
 Operator : JU/SJ
 Sample : J3251-03 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-03-061118

Manual Integrations
 APPROVED

Sohil
 6/14/2018 2:47:45 PM

Quant Time: Jun 14 03:35:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 13 10:51:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	151684	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	585681	20.00	ng	-0.01
38) Acenaphthene-d10	10.01	164	275169	20.00	ng	-0.01
63) Phenanthrene-d10	11.50	188	481685	20.00	ng	0.00
75) Chrysene-d12	14.15	240	387348	20.00	ng	-0.01
86) Perylene-d12	15.69	264	364205	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	417341	46.57	ng	-0.01
7) Phenol-d6	6.61	99	540001	47.69	ng	-0.01
23) Nitrobenzene-d5	7.52	82	328387	32.98	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	126268	47.69	ng	-0.01
44) 2-Fluorobiphenyl	9.32	172	607236	29.27	ng	-0.01
78) Terphenyl-d14	13.08	244	519062	33.28	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	110512	5.621	ng	99
70) Phenanthrene	11.52	178	229246	9.718	ng	97
74) Fluoranthene	12.71	202	180430	7.172	ng	97
77) Pyrene	12.94	202	132702	5.472	ng	99
80) Benzo(a)anthracene	14.14	228	52863	2.293	ng	# 94
82) Chrysene	14.18	228	47435	2.254	ng	# 95
87) Benzo(b)fluoranthene	15.23	252	55248m	2.509	ng	

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

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