

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062118\
 Data File : BF106680.D
 Acq On : 22 Jun 2018 00:02
 Operator : JU/SJ
 Sample : J3245-11 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleID :
 CL-02-061918-B

Manual Integrations
 APPROVED

Sohil
 6/22/2018 10:23:31 AM

Quant Time: Jun 22 04:10:30 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 19 14:32:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	59679	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	219231	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	100309	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	201148	20.00	ng	0.00
75) Chrysene-d12	14.15	240	176941	20.00	ng	0.00
86) Perylene-d12	15.69	264	124747	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	151851	43.09	ng	0.00
7) Phenol-d6	6.60	99	182560	41.48	ng	0.00
23) Nitrobenzene-d5	7.52	82	107831	27.33	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	47744	41.07	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	206953	24.41	ng	-0.01
78) Terphenyl-d14	13.08	244	265738	36.38	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	20584	2.736	ng	Qvalue # 97
70) Phenanthrene	11.52	178	69978	7.213	ng	98
74) Fluoranthene	12.71	202	131684	12.315	ng	96
77) Pyrene	12.94	202	149999	12.856	ng	99
80) Benzo(a)anthracene	14.14	228	61262	5.681	ng	95
82) Chrysene	14.17	228	65892m	6.642	ng	
85) Indeno(1,2,3-cd)pyrene	17.26	276	25845	3.070	ng	# 87
87) Benzo(b)fluoranthene	15.23	252	59521m	7.681	ng	
88) Benzo(k)fluoranthene	15.25	252	18078m	2.450	ng	
89) Benzo(a)pyrene	15.62	252	43067	6.252	ng	# 91
91) Benzo(g,h,i)perylene	17.74	276	31647	5.546	ng	# 89

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

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