

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062118\
 Data File : BF106677.D
 Acq On : 21 Jun 2018 22:42
 Operator : JU/SJ
 Sample : J3523-06 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AREA2(S)

Manual Integrations
 APPROVED

Sohil
 6/22/2018 10:23:24 AM

Quant Time: Jun 22 03:57:45 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	66919	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	257482	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	117675	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	206410	20.00	ng	0.00
75) Chrysene-d12	14.14	240	203364	20.00	ng	-0.01
86) Perylene-d12	15.68	264	149553	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	34019	8.61	ng	0.00
7) Phenol-d6	6.60	99	122855	24.89	ng	-0.01
23) Nitrobenzene-d5	7.52	82	79105	17.07	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	1645	1.21	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	158546	9.75	ng	-0.01
78) Terphenyl-d14	13.07	244	163657	19.49	ng	-0.01
Target Compounds						
70) Phenanthrene	11.52	178	100264	10.071	ng	99
71) Anthracene	11.57	178	27785	2.734	ng	99
74) Fluoranthene	12.71	202	218777	19.938	ng	93
77) Pyrene	12.94	202	204280	15.233	ng	99
80) Benzo(a)anthracene	14.14	228	154718	12.483	ng	99
82) Chrysene	14.17	228	135823	11.911	ng	98
85) Indeno(1,2,3-cd)pyrene	17.24	276	33442	3.456	ng	# 92
87) Benzo(b)fluoranthene	15.22	252	129283m	13.916	ng	
88) Benzo(k)fluoranthene	15.25	252	39969m	4.518	ng	
89) Benzo(a)pyrene	15.61	252	79449	9.620	ng	97
91) Benzo(g,h,i)perylene	17.72	276	31954	4.671	ng	# 88

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

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