

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062118\
 Data File : BF106678.D
 Acq On : 21 Jun 2018 23:09
 Operator : JU/SJ
 Sample : J3523-07 2X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AREA2(D)

Manual Integrations
 APPROVED

Sohil
 6/22/2018 10:23:26 AM

Quant Time: Jun 22 04:02:16 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	66951	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	256119	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	122799	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	220676	20.00	ng	0.00
75) Chrysene-d12	14.15	240	190992	20.00	ng	0.00
86) Perylene-d12	15.69	264	145504	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	95846	24.24	ng	0.00
7) Phenol-d6	6.60	99	121618	24.63	ng	-0.01
23) Nitrobenzene-d5	7.52	82	71628	15.54	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	28257	19.85	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	132303	4.22	ng	-0.01
78) Terphenyl-d14	13.08	244	134996	17.12	ng	0.00
Target Compounds						
31) Naphthalene	8.27	128	34245	2.860	ng	99
49) Dimethylphthalate	9.71	163	21262	2.309	ng	# 95
51) Acenaphthene	10.04	154	43355	5.661	ng	96
54) Dibenzofuran	10.21	168	42264	4.030	ng	94
57) Fluorene	10.55	166	55209	6.634	ng	98
70) Phenanthrene	11.52	178	465142	43.701	ng	98
71) Anthracene	11.57	178	149631	13.773	ng	98
72) Carbazole	11.72	167	31489	3.096	ng	97
74) Fluoranthene	12.72	202	580085	49.447	ng	94
77) Pyrene	12.95	202	584282	46.391	ng	99
80) Benzo(a)anthracene	14.14	228	300089	25.780	ng	99
82) Chrysene	14.18	228	256763	23.976	ng	99
85) Indeno(1,2,3-cd)pyrene	17.25	276	99925	10.995	ng	# 87
87) Benzo(b)fluoranthene	15.23	252	262009m	28.988	ng	
88) Benzo(k)fluoranthene	15.25	252	88101m	10.235	ng	
89) Benzo(a)pyrene	15.62	252	200496	24.952	ng	96
90) Dibenzo(a,h)anthracene	17.25	278	26244m	3.854	ng	
91) Benzo(g,h,i)perylene	17.73	276	100968	15.170	ng	# 88

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

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