

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062618\
 Data File : BF106827.D
 Acq On : 26 Jun 2018 17:14
 Operator : JU/SJ
 Sample : J3684-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleID :
 TP-1-S

Manual Integrations
 APPROVED

Sohil
 6/27/2018 2:14:13 PM

Quant Time: Jun 27 01:41:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 25 18:50:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	47506	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	170637	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	72415	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	127390	20.00	ng	0.00
75) Chrysene-d12	14.16	240	130041	20.00	ng	0.00
86) Perylene-d12	15.71	264	91888	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	275750	98.29	ng	0.00
7) Phenol-d6	6.60	99	340502	97.19	ng	-0.01
23) Nitrobenzene-d5	7.52	82	209054	68.09	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	84594	100.79	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	344129	79.48	ng	-0.01
78) Terphenyl-d14	13.08	244	359879	67.04	ng	0.00
Target Compounds						
48) Acenaphthylene	9.87	152	27970	3.900	ng	98
49) Dimethylphthalate	9.71	163	73276	13.492	ng	# 98
51) Acenaphthene	10.04	154	11923	2.640	ng	98
54) Dibenzofuran	10.21	168	14253	2.305	ng	95
57) Fluorene	10.55	166	18412	3.752	ng	# 98
70) Phenanthrene	11.52	178	224790	36.585	ng	98
71) Anthracene	11.57	178	78164	12.463	ng	98
72) Carbazole	11.72	167	15911	2.710	ng	98
74) Fluoranthene	12.72	202	535798	79.117	ng	96
77) Pyrene	12.95	202	516286	60.206	ng	99
80) Benzo(a)anthracene	14.15	228	348828	44.013	ng	97
82) Chrysene	14.18	228	274369	37.629	ng	97
85) Indeno(1,2,3-cd)pyrene	17.28	276	95844	15.489	ng	# 88
87) Benzo(b)fluoranthene	15.25	252	305307m	53.487	ng	
88) Benzo(k)fluoranthene	15.28	252	86001m	15.821	ng	
89) Benzo(a)pyrene	15.64	252	203234m	40.052	ng	
90) Dibenzo(a,h)anthracene	17.29	278	26418	6.143	ng	# 70
91) Benzo(g,h,i)perylene	17.77	276	90969	21.642	ng	# 89

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

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