

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062218\
 Data File : BF106720.D
 Acq On : 22 Jun 2018 20:36
 Operator : JU/SJ
 Sample : J3246-03 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-062018

Manual Integrations
 APPROVED

Sohil
 6/25/2018 5:16:20 PM

Quant Time: Jun 25 05:44:09 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 22 11:34:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	53076	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	202090	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	93180	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	160774	20.00	ng	0.00
75) Chrysene-d12	14.15	240	145336	20.00	ng	0.00
86) Perylene-d12	15.68	264	122421	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	144478	46.09	ng	0.00
7) Phenol-d6	6.60	99	183728	46.94	ng	-0.01
23) Nitrobenzene-d5	7.52	82	112531	30.95	ng	-0.01
41) 2,4,6-Tribromophenol	10.79	330	47049	43.56	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	226229	31.88	ng	-0.01
78) Terphenyl-d14	13.08	244	223703	37.28	ng	0.00
Target Compounds						
31) Naphthalene	8.26	128	77768	8.231	ng	99
37) 2-Methylnaphthalene	8.95	142	54459	8.696	ng	94
48) Acenaphthylene	9.86	152	23928	2.593	ng	# 96
49) Dimethylphthalate	9.71	163	20920	2.993	ng	# 97
51) Acenaphthene	10.03	154	74929	12.894	ng	97
54) Dibenzofuran	10.20	168	71586	8.996	ng	94
57) Fluorene	10.55	166	113512	17.976	ng	99
70) Phenanthrene	11.52	178	703009	90.659	ng	99
71) Anthracene	11.57	178	193490	24.446	ng	98
72) Carbazole	11.72	167	115215	15.548	ng	97
74) Fluoranthene	12.72	202	771941	90.318	ng	95
77) Pyrene	12.95	202	707043	73.774	ng	98
80) Benzo(a)anthracene	14.14	228	374807	42.314	ng	97
82) Chrysene	14.18	228	327262	40.159	ng	97
85) Indeno(1,2,3-cd)pyrene	17.25	276	136698	19.767	ng	# 87
87) Benzo(b)fluoranthene	15.23	252	366463m	48.189	ng	
88) Benzo(k)fluoranthene	15.26	252	133467m	18.430	ng	
89) Benzo(a)pyrene	15.62	252	254480m	37.643	ng	
90) Dibenzo(a,h)anthracene	17.26	278	33195	5.794	ng	# 91
91) Benzo(a,h,i)perylene	17.74	276	132318	23.628	ng	# 89

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

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