

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\
 Data File : BF107472.D
 Acq On : 18 Jul 2018 12:54
 Operator : JU/SJ
 Sample : J4016-11 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB3-L-12-7

Manual Integrations
 APPROVED

Sohil
 7/19/2018 11:49:25 AM

Quant Time: Jul 18 14:38:26 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 17 11:29:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	139573	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	525979	20.00	ng	0.00
38) Acenaphthene-d10	10.03	164	299085	20.00	ng	0.00
63) Phenanthrene-d10	11.52	188	594838	20.00	ng	0.00
75) Chrysene-d12	14.18	240	328068	20.00	ng	0.00
86) Perylene-d12	15.72	264	344584	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	360329	39.21	ng	0.00
7) Phenol-d6	6.61	99	473335	39.84	ng	0.00
23) Nitrobenzene-d5	7.54	82	336586	34.12	ng	-0.01
41) 2,4,6-Tribromophenol	10.82	330	92413	35.17	ng	0.00
44) 2-Fluorobiphenyl	9.34	172	544716	22.38	ng	0.00
78) Terphenyl-d14	13.11	244	521272	39.48	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.29	128	164524	6.440	ng	96
37) 2-Methylnaphthalene	8.98	142	50556	3.002	ng	# 89
48) Acenaphthylene	9.89	152	116200	4.077	ng	97
49) Dimethylphthalate	9.73	163	104825	4.411	ng	# 94
51) Acenaphthene	10.06	154	106149	5.618	ng	92
54) Dibenzofuran	10.23	168	160750	5.505	ng	91
57) Fluorene	10.58	166	191007	9.877	ng	92
70) Phenanthrene	11.56	178	2088854	70.464	ng	98
71) Anthracene	11.60	178	569397	19.284	ng	97
72) Carbazole	11.75	167	179373	6.417	ng	100
74) Fluoranthene	12.75	202	2482183	73.670	ng	93
77) Pyrene	12.98	202	2084490	93.436	ng	97
80) Benzo(a)anthracene	14.17	228	1052590	52.438	ng	95
82) Chrysene	14.21	228	900224	46.948	ng	98
85) Indeno(1,2,3-cd)pyrene	17.31	276	510271	37.386	ng	# 82
87) Benzo(b)fluoranthene	15.27	252	1061274m	52.619	ng	
88) Benzo(k)fluoranthene	15.29	252	431573m	21.926	ng	
89) Benzo(a)pyrene	15.66	252	837177m	45.223	ng	
90) Dibenzo(a,h)anthracene	17.31	278	118449	8.478	ng	96
91) Benzo(a,h,i)perylene	17.79	276	444725	31.327	ng	# 88

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

