

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032919\
 Data File : BF113449.D
 Acq On : 30 Mar 2019 9:13
 Operator : JU/SJ
 Sample : K2153-01
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CONC-1

Manual Integrations
 APPROVED

Sohil
 4/1/2019 5:46:43 PM

Quant Time: Apr 01 04:40:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	100985	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	384235	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	184898	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	351240	20.00	ng	0.00
76) Chrysene-d12	13.85	240	310961	20.00	ng	0.00
87) Perylene-d12	15.25	264	261030	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	694287	112.42	ng	0.00
7) Phenol-d6	6.35	99	960085	122.88	ng	0.00
23) Nitrobenzene-d5	7.27	82	580972	89.74	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	193049	84.53	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	1111610	96.65	ng	0.00
79) Terphenyl-d14	12.79	244	1170670	82.98	ng	0.00

Target Compounds				Qvalue		
31) Naphthalene	8.00	128	74769	3.982	ng	99
37) 2-Methylnaphthalene	8.70	142	39741	3.392	ng	97
38) 1-Methylnaphthalene	8.80	142	43574	3.867	ng	98
49) Acenaphthylene	9.60	152	73327	3.814	ng	98
50) Dimethylphthalate	9.45	163	90412	6.598	ng	100
52) Acenaphthene	9.76	154	91575	8.458	ng	99
55) Dibenzofuran	9.94	168	68974	4.238	ng	# 94
58) Fluorene	10.28	166	102633	8.117	ng	98
71) Phenanthrene	11.24	178	1527313	87.562	ng	99
72) Anthracene	11.29	178	385804	21.776	ng	99
73) Carbazole	11.45	167	132551	7.841	ng	99
75) Fluoranthene	12.43	202	2156384	109.358	ng	98
78) Pvrene	12.66	202	2230482	103.510	ng	99
80) Butylbenzylphthalate	13.27	149	65512	6.898	ng	95
81) Benzo(a)anthracene	13.84	228	1209159	67.931	ng	97
83) Chrysene	13.87	228	1079183	59.690	ng	97
84) Bis(2-ethylhexyl)phthalate	13.83	149	116444	10.002	ng	100
86) Indeno(1,2,3-cd)pyrene	16.62	276	570671	36.778	ng	95
88) Benzo(b)fluoranthene	14.86	252	1260347m	77.727	ng	
89) Benzo(k)fluoranthene	14.88	252	453033m	31.815	ng	
90) Benzo(a)pyrene	15.20	252	915521	63.687	ng	98
91) Dibenzo(a,h)anthracene	16.62	278	141028m	11.346	ng	
92) Benzo(g,h,i)perylene	17.03	276	558433	44.558	ng	97

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032919\
Data File : BF113449.D
Acq On : 30 Mar 2019 9:13
Operator : JU/SJ
Sample : K2153-01
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CONC-1

Manual Integrations
APPROVED

Sohil
4/1/2019 5:46:43 PM

Quant Time: Apr 01 04:40:20 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF032519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 26 03:32:08 2019
Response via : Initial Calibration

