

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060119\
 Data File : BF114763.D
 Acq On : 1 Jun 2019 17:59
 Operator : HP/JU
 Sample : K3105-05
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRIANGLE

Manual Integrations
 APPROVED

mohammad
 6/3/2019 8:51:38 AM

Quant Time: Jun 03 04:20:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 19:08:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	339609	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	1175799	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	497402	20.00	ng	0.00
64) Phenanthrene-d10	11.20	188	935911	20.00	ng	0.00
76) Chrysene-d12	13.83	240	674147	20.00	ng	0.02
87) Perylene-d12	15.22	264	666298	20.00	ng	0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	1806365	93.39	ng	0.01
7) Phenol-d6	6.33	99	2126833	91.24	ng	0.00
23) Nitrobenzene-d5	7.25	82	1257789	66.74	ng	0.00
42) 2,4,6-Tribromophenol	10.50	330	462284	97.54	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	1921826	69.70	ng	0.00
79) Terphenyl-d14	12.78	244	1839533	51.30	ng	0.00

Target Compounds				Qvalue		
31) Naphthalene	8.00	128	345564	6.381	ng	96
37) 2-Methylnaphthalene	8.69	142	536594	14.912	ng	98
38) 1-Methylnaphthalene	8.79	142	750444	22.011	ng	99
46) 1,1'-Biphenyl	9.15	154	126726	3.432	ng	96
49) Acenaphthylene	9.58	152	768021	16.414	ng	96
50) Dimethylphthalate	9.44	163	323972	9.347	ng	100
52) Acenaphthene	9.75	154	568329	19.806	ng	98
55) Dibenzofuran	9.92	168	331252	8.110	ng	# 92
58) Fluorene	10.27	166	1045118	36.754	ng	99
71) Phenanthrene	11.23	178	5250119	109.968	ng	94
72) Anthracene	11.27	178	2035897	42.134	ng	99
73) Carbazole	11.42	167	236778	5.576	ng	99
75) Fluoranthene	12.42	202	5240226	106.000	ng	96
78) Pyrene	12.65	202	6001257	105.234	ng	93
80) Butylbenzylphthalate	13.26	149	107592	3.754	ng	94
81) Benzo(a)anthracene	13.82	228	3771617	72.803	ng	92
83) Chrysene	13.86	228	3370804	68.102	ng	96
86) Indeno(1,2,3-cd)pyrene	16.55	276	1184965	23.903	ng	94
88) Benzo(b)fluoranthene	14.84	252	3164766m	75.413	ng	
89) Benzo(k)fluoranthene	14.86	252	1039701m	28.504	ng	
90) Benzo(a)pyrene	15.17	252	2357852	63.956	ng	96
91) Dibenz(a,h)anthracene	16.56	278	356996	10.958	ng	94
92) Benzo(g,h,i)perylene	16.95	276	1129442	35.721	ng	99

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060119\
Data File : BF114763.D
Acq On : 1 Jun 2019 17:59
Operator : HP/JU
Sample : K3105-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TRIANGLE

Quant Time: Jun 03 04:20:30 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 29 19:08:51 2019
Response via : Initial Calibration

Manual Integrations
APPROVED

mohammad
6/3/2019 8:51:38 AM

