

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112619\
 Data File : BF118018.D
 Acq On : 26 Nov 2019 20:20
 Operator : JU
 Sample : K5938-01 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ETGI-277

Manual Integrations
 APPROVED

jagrut
 11/27/2019 11:33:49 PM

Quant Time: Nov 27 01:17:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 26 05:38:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	123773	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	461609	20.00	ng	0.00
39) Acenaphthene-d10	9.95	164	247266	20.00	ng	0.00
64) Phenanthrene-d10	11.44	188	372746	20.00	ng	0.00
76) Chrysene-d12	14.10	240	254527	20.00	ng	0.01
87) Perylene-d12	15.60	264	190670	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	316823	45.31	ng	0.00
7) Phenol-d6	6.52	99	394665	46.79	ng	0.00
23) Nitrobenzene-d5	7.45	82	228965	29.49	ng	-0.01
42) 2,4,6-Tribromophenol	10.74	330	87187	33.31	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	507694	36.84	ng	0.00
79) Terphenyl-d14	13.03	244	358832	25.77	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.20	128	52745	2.451	ng	98
37) 2-Methylnaphthalene	8.90	142	57918	3.983	ng	97
38) 1-Methylnaphthalene	9.00	142	84615	6.155	ng	99
49) Acenaphthylene	9.80	152	454202	20.643	ng	98
50) Dimethylphthalate	9.65	163	71509	4.124	ng	# 98
52) Acenaphthene	9.98	154	40212	2.853	ng	94
58) Fluorene	10.50	166	126016	8.332	ng	98
71) Phenanthrene	11.47	178	1175081	62.703	ng	98
72) Anthracene	11.52	178	303920	16.357	ng	98
73) Carbazole	11.67	167	34065	2.098	ng	93
75) Fluoranthene	12.67	202	1032607	63.622	ng	98
78) Pyrene	12.90	202	1494391	72.650	ng	99
81) Benzo(a)anthracene	14.09	228	574196m	34.139	ng	
83) Chrysene	14.13	228	694598	42.618	ng	98
86) Indeno(1,2,3-cd)pyrene	17.14	276	192540	13.880	ng	95
88) Benzo(b)fluoranthene	15.16	252	450348m	36.354	ng	
89) Benzo(k)fluoranthene	15.19	252	124161m	10.637	ng	
90) Benzo(a)pyrene	15.54	252	378605	34.756	ng	98
91) Dibenzo(a,h)anthracene	17.14	278	56702	6.048	ng	# 81
92) Benzo(a,h,i)perylene	17.60	276	213282	22.838	ng	98

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112619\
Data File : BF118018.D
Acq On : 26 Nov 2019 20:20
Operator : JU
Sample : K5938-01 2X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-277

Manual Integrations
APPROVED

jagrut
11/27/2019 11:33:49 PM

Quant Time: Nov 27 01:17:30 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 26 05:38:20 2019
Response via : Initial Calibration

