

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122619\
 Data File : BF118318.D
 Acq On : 26 Dec 2019 16:03
 Operator : JU
 Sample : K6395-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1

Manual Integrations
 APPROVED

mohammad
 12/27/2019 3:19:12 PM

Quant Time: Dec 27 04:57:16 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 15:58:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	119203	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	485977	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	262996	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	503592	20.00	ng	0.00
76) Chrysene-d12	14.04	240	333112	20.00	ng	-0.01
87) Perylene-d12	15.53	264	357892	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	704857	100.94	ng	0.00
7) Phenol-d6	6.48	99	902126	104.14	ng	0.00
23) Nitrobenzene-d5	7.41	82	535486	63.02	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	323941	99.74	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	1165325	67.64	ng	0.00
79) Terphenyl-d14	12.98	244	1444176	71.93	ng	0.00
Target Compounds						
10) Phenol	6.49	94	31095	3.194	ng	97
50) Dimethylphthalate	9.61	163	135230	6.850	ng	100
71) Phenanthrene	11.41	178	553311	20.155	ng	99
72) Anthracene	11.46	178	116690	4.247	ng	99
75) Fluoranthene	12.61	202	622093	22.453	ng	99
78) Pyrene	12.84	202	623481	22.920	ng	99
81) Benzo(a)anthracene	14.03	228	270040	11.581	ng	100
83) Chrysene	14.07	228	244999	10.974	ng	99
86) Indeno(1,2,3-cd)pyrene	17.03	276	123721	6.504	ng	95
88) Benzo(b)fluoranthene	15.10	252	278399m	11.719	ng	
89) Benzo(k)fluoranthene	15.12	252	72585m	3.182	ng	
90) Benzo(a)pyrene	15.47	252	209505	10.011	ng	95
91) Dibenzo(a,h)anthracene	17.03	278	36524m	2.011	ng	
92) Benzo(g,h,i)perylene	17.47	276	123095	7.017	ng	97

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

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