

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082319\
 Data File : BF116497.D
 Acq On : 24 Aug 2019 00:25
 Operator : HP/JU
 Sample : K4403-16
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 T1-J-(0-2.5)

Manual Integrations
 APPROVED

Jagrut
 8/28/2019 8:10:49 AM

Quant Time: Aug 24 04:45:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 23 13:38:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	142246	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	555276	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	270003	20.00	ng	0.00
64) Phenanthrene-d10	11.40	188	415102	20.00	ng	0.00
76) Chrysene-d12	14.03	240	234557	20.00	ng	0.00
87) Perylene-d12	15.50	264	234798	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	671067	68.93	ng	0.01
7) Phenol-d6	6.51	99	903631	75.70	ng	0.00
23) Nitrobenzene-d5	7.44	82	551762	54.36	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	156298	67.42	ng	0.00
45) 2-Fluorobiphenyl	9.24	172	769584	45.90	ng	0.00
79) Terphenyl-d14	12.98	244	658768	52.46	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.63	163	126506	6.715	ng	99
71) Phenanthrene	11.42	178	407283	18.011	ng	99
72) Anthracene	11.47	178	92733	4.081	ng	99
75) Fluoranthene	12.61	202	557767	23.858	ng	98
78) Pyrene	12.84	202	543584	28.979	ng	99
81) Benzo(a)anthracene	14.02	228	241283	15.883	ng	98
83) Chrysene	14.06	228	197469	12.825	ng	97
86) Indeno(1,2,3-cd)pyrene	16.96	276	89183	6.762	ng	95
88) Benzo(b)fluoranthene	15.07	252	220842m	15.226	ng	
89) Benzo(k)fluoranthene	15.10	252	60353m	4.593	ng	
90) Benzo(a)pyrene	15.44	252	147765	11.527	ng	# 95
91) Dibenzo(a,h)anthracene	16.97	278	23575	2.238	ng	# 65
92) Benzo(a,h,i)perylene	17.40	276	83002	8.265	ng	99

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

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