

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123019\
 Data File : BF118391.D
 Acq On : 30 Dec 2019 20:01
 Operator : JU
 Sample : K6455-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 K435-WC-1

Manual Integrations
 APPROVED

mohammad
 12/31/2019 12:04:29 PM

Quant Time: Dec 31 02:37:24 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	138346	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	538890	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	282014	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	487132	20.00	ng	0.00
76) Chrysene-d12	14.05	240	286985	20.00	ng	0.00
87) Perylene-d12	15.54	264	325816	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	835364	103.07	ng	0.01
7) Phenol-d6	6.49	99	1094514	108.86	ng	0.00
23) Nitrobenzene-d5	7.42	82	676289	71.78	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	300996	86.43	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	1415844	76.64	ng	0.00
79) Terphenyl-d14	12.99	244	1353560	78.25	ng	0.00
Target Compounds						
15) Benzyl Alcohol	7.00	79	195542	25.544	ng	98
50) Dimethylphthalate	9.61	163	99048	4.679	ng	98
71) Phenanthrene	11.42	178	157653	5.937	ng	98
75) Fluoranthene	12.62	202	244359	9.118	ng	99
78) Pyrene	12.84	202	241210	10.293	ng	99
81) Benzo(a)anthracene	14.04	228	112202	5.586	ng	97
83) Chrysene	14.07	228	100041	5.201	ng	99
86) Indeno(1,2,3-cd)pyrene	17.06	276	46728	2.851	ng	95
88) Benzo(b)fluoranthene	15.10	252	136948m	6.332	ng	
90) Benzo(a)pyrene	15.48	252	90517	4.751	ng	95
92) Benzo(g,h,i)perylene	17.52	276	43929	2.751	ng	95

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

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