

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010220\
 Data File : BF118445.D
 Acq On : 2 Jan 2020 19:47
 Operator : JU
 Sample : K6465-06
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GP-9-(8-10)

Manual Integrations
 APPROVED

mohammad
 1/3/2020 11:04:25 AM

Quant Time: Jan 03 03:32:57 2020
 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	125422	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	488307	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	249861	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	421206	20.00	ng	0.00
76) Chrysene-d12	14.04	240	277868	20.00	ng	0.00
87) Perylene-d12	15.53	264	280196	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	644505	87.72	ng	0.00
7) Phenol-d6	6.49	99	830346	91.10	ng	0.00
23) Nitrobenzene-d5	7.41	82	479137	56.12	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	233798	75.77	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	939828	57.42	ng	-0.01
79) Terphenyl-d14	12.97	244	850114	50.76	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.61	163	60604	3.231	ng	99
71) Phenanthrene	11.41	178	82829	3.607	ng	98
75) Fluoranthene	12.61	202	151574	6.541	ng	96
78) Pyrene	12.84	202	125233	5.519	ng	98
81) Benzo(a)anthracene	14.03	228	56814	2.921	ng	96
83) Chrysene	14.07	228	47375	2.544	ng	97
88) Benzo(b)fluoranthene	15.10	252	62430m	3.357	ng	
90) Benzo(a)pyrene	15.47	252	46566	2.842	ng	95

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

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