

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010220\
 Data File : BF118451.D
 Acq On : 2 Jan 2020 22:34
 Operator : JU
 Sample : K6465-05
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GP-9-(1-3)

Manual Integrations
 APPROVED

mohammad
 1/3/2020 11:04:46 AM

Quant Time: Jan 03 04:02:04 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	83355	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	326173	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	172279	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	285878	20.00	ng	0.00
76) Chrysene-d12	14.04	240	190088	20.00	ng	-0.01
87) Perylene-d12	15.53	264	200138	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	487402	99.81	ng	0.00
7) Phenol-d6	6.49	99	600918	99.20	ng	0.00
23) Nitrobenzene-d5	7.42	82	361563	63.40	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	174269	81.91	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	721162	63.90	ng	-0.01
79) Terphenyl-d14	12.97	244	688588	60.10	ng	-0.01
Target Compounds						
10) Phenol	6.50	94	20894	3.069	ng	85
50) Dimethylphthalate	9.61	163	80369	6.215	ng	100
71) Phenanthrene	11.41	178	57824	3.710	ng	98
75) Fluoranthene	12.61	202	97983	6.230	ng	99
78) Pyrene	12.84	202	83468	5.377	ng	99
81) Benzo(a)anthracene	14.03	228	41292	3.103	ng	97
83) Chrysene	14.07	228	40962	3.215	ng	98
88) Benzo(b)fluoranthene	15.09	252	55022m	4.142	ng	
90) Benzo(a)pyrene	15.47	252	34710	2.966	ng	96

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

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