

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010220\
 Data File : BF118448.D
 Acq On : 2 Jan 2020 21:11
 Operator : JU
 Sample : K6465-07
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GP-4-(1-3)

Manual Integrations
 APPROVED

mohammad
 1/3/2020 11:04:33 AM

Quant Time: Jan 03 03:45:48 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	96883	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	360661	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	204375	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	331654	20.00	ng	0.00
76) Chrysene-d12	14.04	240	214485	20.00	ng	0.00
87) Perylene-d12	15.53	264	205885	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	531241	93.60	ng	0.00
7) Phenol-d6	6.49	99	698007	99.14	ng	0.00
23) Nitrobenzene-d5	7.41	82	430467	68.27	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	199423	79.01	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	717789	53.61	ng	-0.01
79) Terphenyl-d14	12.97	244	674812	52.20	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.61	163	63218	4.121	ng	98
71) Phenanthrene	11.41	178	95666	5.291	ng	98
75) Fluoranthene	12.61	202	100632	5.515	ng	100
78) Pyrene	12.84	202	92772	5.297	ng	99
81) Benzo(a)anthracene	14.03	228	43946	2.927	ng	98
83) Chrysene	14.07	228	41592m	2.893	ng	
88) Benzo(b)fluoranthene	15.10	252	50108m	3.666	ng	
90) Benzo(a)pyrene	15.47	252	31184	2.590	ng	97
92) Benzo(g,h,i)perylene	17.50	276	20572	2.038	ng	# 91

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

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