

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122619\
 Data File : BF118324.D
 Acq On : 26 Dec 2019 18:51
 Operator : JU
 Sample : K6437-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GP-5-(5-6)

Manual Integrations
 APPROVED

mohammad
 12/27/2019 3:19:22 PM

Quant Time: Jan 09 06:57:46 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	97725	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	378668	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	200768	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	354038	20.00	ng	0.00
76) Chrysene-d12	14.04	240	267989	20.00	ng	0.00
87) Perylene-d12	15.54	264	282528	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	556029	97.12	ng	0.00
7) Phenol-d6	6.48	99	696913	98.13	ng	0.00
23) Nitrobenzene-d5	7.41	82	394229	59.55	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	215714	87.00	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	794599	60.41	ng	0.00
79) Terphenyl-d14	12.98	244	852768	52.79	ng	0.00
Target Compounds						
10) Phenol	6.49	94	20937	2.623	ng	93
50) Dimethylphthalate	9.61	163	51267	3.402	ng	# 80
75) Fluoranthene	12.61	202	151476	7.777	ng	98
78) Pyrene	12.84	202	339685	15.522	ng	99
81) Benzo(a)anthracene	14.03	228	42688	2.276	ng	# 79
83) Chrysene	14.07	228	45451	2.530	ng	# 83
88) Benzo(b)fluoranthene	15.10	252	63275m	3.374	ng	
90) Benzo(a)pyrene	15.47	252	39083	2.366	ng	# 57
92) Benzo(g,h,i)perylene	17.50	276	41551	3.000	ng	# 92

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

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