

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012420\
 Data File : BF118699.D
 Acq On : 24 Jan 2020 21:35
 Operator : CG/JU
 Sample : L1242-12
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20200065-AMENDED-COMP-3

Manual Integrations
 APPROVED

mohammad
 1/28/2020 8:12:40 AM

Quant Time: Jan 27 02:54:29 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 20 14:45:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	251447	20.00	ng	-0.02
21) Naphthalene-d8	8.15	136	945723	20.00	ng	-0.02
39) Acenaphthene-d10	9.90	164	517522	20.00	ng	-0.03
64) Phenanthrene-d10	11.38	188	890115	20.00	ng	-0.03
76) Chrysene-d12	14.02	240	620636	20.00	ng	-0.03
87) Perylene-d12	15.48	264	697100	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	62559	4.62	ng	-0.02
7) Phenol-d6	6.49	99	560743	31.87	ng	-0.02
23) Nitrobenzene-d5	7.42	82	994696	58.83	ng	-0.02
42) 2,4,6-Tribromophenol	10.68	330	13192	2.20	ng	-0.03
45) 2-Fluorobiphenyl	9.22	172	1651826	52.46	ng	-0.02
79) Terphenyl-d14	12.96	244	1844712	64.83	ng	-0.03
Target Compounds						
10) Phenol	6.50	94	72252	3.603	ng	95
20) 3+4-Methylphenols	7.26	107	88352	5.717	ng	# 57
50) Dimethylphthalate	9.61	163	351819	10.434	ng	97
75) Fluoranthene	12.59	202	200994	4.438	ng	94
78) Pyrene	12.82	202	176923	4.199	ng	# 91
81) Benzo(a)anthracene	14.00	228	97833	2.620	ng	95
83) Chrysene	14.04	228	89450	2.500	ng	93
88) Benzo(b)fluoranthene	15.05	252	100639m	2.297	ng	

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

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