

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111920\
 Data File : BF122396.D
 Acq On : 19 Nov 2020 23:15
 Operator : JU/CG
 Sample : L4802-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-064-B

Manual Integrations
 APPROVED

mohammad
 11/24/2020 8:54:42 AM

Quant Time: Nov 20 00:05:47 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 10 12:25:42 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	203622	20.00	ng	0.00
21) Naphthalene-d8	8.151	136	758763	20.00	ng	0.00
39) Acenaphthene-d10	9.910	164	393003	20.00	ng	0.00
64) Phenanthrene-d10	11.392	188	661296	20.00	ng	0.00
76) Chrysene-d12	14.033	240	603485	20.00	ng	0.00
86) Perylene-d12	15.509	264	567765	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	1045551	78.84	ng	0.02
7) Phenol-d6	6.510	99	1317358	75.94	ng	0.02
23) Nitrobenzene-d5	7.428	82	814370	65.71	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	347720	82.44	ng	0.00
45) 2-Fluorobiphenyl	9.227	172	1486035	59.08	ng	0.00
79) Terphenyl-d14	12.980	244	1392941	48.90	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.522	94	64116	3.75	ng	# 74
50) Dimethylphthalate	9.621	163	85541	3.06	ng	98
75) Fluoranthene	12.604	202	178858	4.57	ng	98
78) Pyrene	12.833	202	175378	4.14	ng	99
81) Benzo(a)anthracene	14.021	228	109953	2.93	ng	94
83) Chrysene	14.057	228	98312	2.60	ng	96
88) Benzo(b)fluoranthene	15.074	252	150816m	4.10	ng	
90) Benzo(a)pyrene	15.445	252	100623	3.14	ng	# 83

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

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