

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
 Data File : BF117734.D
 Acq On : 8 Nov 2019 16:43
 Operator : JU
 Sample : K5722-14
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 2-A

Manual Integrations
 APPROVED

mohammad
 11/11/2019 8:06:04 AM

Quant Time: Nov 09 02:30:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 17:27:10 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	348004	20.00	ng	-0.01
21) Naphthalene-d8	8.22	136	1397902	20.00	ng	-0.01
39) Acenaphthene-d10	9.97	164	764403	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	1318094	20.00	ng	0.00
76) Chrysene-d12	14.12	240	862303	20.00	ng	0.00
87) Perylene-d12	15.62	264	833963	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	1982089	103.10	ng	0.00
7) Phenol-d6	6.55	99	2560520	108.02	ng	-0.01
23) Nitrobenzene-d5	7.49	82	1591114	67.20	ng	-0.02
42) 2,4,6-Tribromophenol	10.77	330	774657	103.60	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	2769544	71.49	ng	-0.01
79) Terphenyl-d14	13.06	244	2713348	64.78	ng	0.00
Target Compounds						
10) Phenol	6.56	94	89317	3.497	ng	Qvalue 96
50) Dimethylphthalate	9.69	163	360568	6.977	ng	96
75) Fluoranthene	12.69	202	311489	4.828	ng	96
78) Pyrene	12.92	202	313221	4.976	ng	96
81) Benzo(a)anthracene	14.11	228	133978	2.496	ng	93
83) Chrysene	14.14	228	159742	3.014	ng	96
88) Benzo(b)fluoranthene	15.17	252	199094m	3.791	ng	
90) Benzo(a)pyrene	15.56	252	91193	2.003	ng	# 96

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

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