

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090119\
 Data File : BF116741.D
 Acq On : 1 Sep 2019 14:38
 Operator : HP/JU
 Sample : K4525-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 T9-12-13-A1-A4-(0-1.9)

Manual Integrations
 APPROVED

mohammad
 9/5/2019 9:54:17 AM

Quant Time: Sep 02 05:38:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 30 20:02:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	120370	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	486072	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	256017	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	435843	20.00	ng	-0.01
76) Chrysene-d12	13.99	240	352230	20.00	ng	-0.01
87) Perylene-d12	15.44	264	324415	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	752215	101.24	ng	0.00
7) Phenol-d6	6.47	99	1048449	107.46	ng	0.00
23) Nitrobenzene-d5	7.40	82	650579	76.41	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	209901	122.64	ng	-0.01
45) 2-Fluorobiphenyl	9.20	172	902078	59.44	ng	-0.01
79) Terphenyl-d14	12.94	244	946142	49.69	ng	-0.01
Target Compounds						
10) Phenol	6.48	94	49448	4.577	ng	98
50) Dimethylphthalate	9.59	163	93719	5.302	ng	100
71) Phenanthrene	11.38	178	67469	2.781	ng	98
75) Fluoranthene	12.56	202	146213	6.132	ng	96
78) Pyrene	12.79	202	136320	4.354	ng	99
81) Benzo(a)anthracene	13.97	228	60594	2.718	ng	99
83) Chrysene	14.01	228	64413m	2.805	ng	
84) Bis(2-ethylhexyl)phthalate	13.97	149	102266	5.441	ng	99
88) Benzo(b)fluoranthene	15.02	252	74638m	3.805	ng	
90) Benzo(a)pyrene	15.37	252	48192	2.824	ng	# 90
92) Benzo(g,h,i)perylene	17.30	276	33639	2.208	ng	# 88

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

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