

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111219\
 Data File : BF117764.D
 Acq On : 12 Nov 2019 17:22
 Operator : JU
 Sample : K5788-06
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-1

Manual Integrations
 APPROVED

mohammad
 11/13/2019 6:48:31 PM

Quant Time: Nov 13 03:42:06 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 12 13:32:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	278266	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	1043971	20.00	ng	0.00
39) Acenaphthene-d10	9.97	164	509056	20.00	ng	0.00
64) Phenanthrene-d10	11.46	188	829343	20.00	ng	0.00
76) Chrysene-d12	14.12	240	634758	20.00	ng	0.00
87) Perylene-d12	15.63	264	756226	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	1374350	89.40	ng	0.00
7) Phenol-d6	6.55	99	1696707	89.51	ng	0.00
23) Nitrobenzene-d5	7.49	82	1049360	59.34	ng	0.00
42) 2,4,6-Tribromophenol	10.76	330	420501	84.44	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	1924009	74.57	ng	0.00
79) Terphenyl-d14	13.06	244	1776823	57.62	ng	0.00
Target Compounds						
10) Phenol	6.55	94	89021	4.359	ng	Qvalue 85
50) Dimethylphthalate	9.68	163	239371	6.956	ng	97
71) Phenanthrene	11.49	178	84770	2.132	ng	96
75) Fluoranthene	12.68	202	261481	6.442	ng	97
78) Pyrene	12.91	202	238488	5.147	ng	95
81) Benzo(a)anthracene	14.10	228	155900	3.946	ng	98
83) Chrysene	14.14	228	148262	3.800	ng	97
86) Indeno(1,2,3-cd)pyrene	17.15	276	92080	2.535	ng	95
88) Benzo(b)fluoranthene	15.17	252	199064m	4.180	ng	
90) Benzo(a)pyrene	15.56	252	142307	3.447	ng	# 95
92) Benzo(g,h,i)perylene	17.62	276	84419	2.352	ng	96

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

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