

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101219\
 Data File : BF117386.D
 Acq On : 12 Oct 2019 23:17
 Operator : JU
 Sample : K5316-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-9

Manual Integrations
 APPROVED

mohammad
 10/16/2019 8:06:02 AM

Quant Time: Oct 14 03:22:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 09 01:26:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	39127	20.00	ng	-0.01
21) Naphthalene-d8	8.07	136	160702	20.00	ng	-0.01
39) Acenaphthene-d10	9.83	164	88078	20.00	ng	-0.01
64) Phenanthrene-d10	11.32	188	151434	20.00	ng	-0.01
76) Chrysene-d12	13.96	240	95640	20.00	ng	-0.01
87) Perylene-d12	15.42	264	97566	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	240620	96.01	ng	-0.01
7) Phenol-d6	6.44	99	341400	105.41	ng	-0.01
23) Nitrobenzene-d5	7.36	82	212148	69.36	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	70809	96.19	ng	-0.01
45) 2-Fluorobiphenyl	9.15	172	328706	58.35	ng	-0.01
79) Terphenyl-d14	12.90	244	289111	56.50	ng	-0.01
Target Compounds						
10) Phenol	6.45	94	8801	2.465	ng	# 78
49) Acenaphthylene	9.69	152	18875	2.311	ng	95
50) Dimethylphthalate	9.55	163	61810	9.911	ng	# 98
71) Phenanthrene	11.34	178	88873	10.332	ng	98
75) Fluoranthene	12.53	202	106586	12.060	ng	98
78) Pyrene	12.76	202	142576	17.767	ng	98
81) Benzo(a)anthracene	13.95	228	51574	8.071	ng	99
83) Chrysene	13.99	228	48331	7.755	ng	100
86) Indeno(1,2,3-cd)pyrene	16.87	276	20624	4.536	ng	96
88) Benzo(b)fluoranthene	15.00	252	37462m	6.339	ng	
89) Benzo(k)fluoranthene	15.02	252	12011m	2.145	ng	
90) Benzo(a)pyrene	15.36	252	35064	6.761	ng	# 96
92) Benzo(a,h,i)perylene	17.31	276	21884	5.316	ng	96

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

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