

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111219\
 Data File : BF117765.D
 Acq On : 12 Nov 2019 17:49
 Operator : JU
 Sample : K5791-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FINES-NORTH-BERGEN

Manual Integrations
 APPROVED

mohammad
 11/13/2019 6:48:34 PM

Quant Time: Nov 13 01:34:24 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	276799	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	1054870	20.00	ng	0.00
39) Acenaphthene-d10	9.97	164	526054	20.00	ng	0.00
64) Phenanthrene-d10	11.46	188	842583	20.00	ng	0.00
76) Chrysene-d12	14.12	240	597461	20.00	ng	0.00
87) Perylene-d12	15.63	264	697288	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	1330987	87.04	ng	0.00
7) Phenol-d6	6.55	99	1665283	88.32	ng	0.00
23) Nitrobenzene-d5	7.49	82	1024348	57.33	ng	0.00
42) 2,4,6-Tribromophenol	10.76	330	419684	81.56	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	1861563	69.82	ng	0.00
79) Terphenyl-d14	13.06	244	1620880	55.85	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.68	163	211051	5.935	ng	98
71) Phenanthrene	11.49	178	967743	23.962	ng	97
72) Anthracene	11.54	178	130855	3.257	ng	96
75) Fluoranthene	12.69	202	1394983	33.827	ng	97
78) Pyrene	12.92	202	1192091	27.333	ng	96
81) Benzo(a)anthracene	14.10	228	572179	15.385	ng	98
83) Chrysene	14.14	228	560842	15.271	ng	97
86) Indeno(1,2,3-cd)pyrene	17.16	276	279618	8.180	ng	94
88) Benzo(b)fluoranthene	15.18	252	758469m	17.271	ng	
89) Benzo(k)fluoranthene	15.20	252	251672m	6.425	ng	
90) Benzo(a)pyrene	15.56	252	490150	12.875	ng	99
91) Dibenzo(a,h)anthracene	17.17	278	72082m	2.111	ng	
92) Benzo(a,h,i)perylene	17.62	276	234928	7.100	ng	98

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

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