

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010919\
 Data File : BF111941.D
 Acq On : 9 Jan 2019 10:45
 Operator : JU/SJ
 Sample : K1066-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TN-1419

Manual Integrations
 APPROVED

Sohil
 1/10/2019 11:45:42 AM

Quant Time: Jan 09 11:20:45 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 07 14:33:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	200018	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	845417	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	459827	20.00	ng	-0.01
64) Phenanthrene-d10	11.41	188	865039	20.00	ng	0.00
76) Chrysene-d12	14.05	240	578764	20.00	ng	0.00
87) Perylene-d12	15.54	264	467990	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	1079611	92.34	ng	0.00
7) Phenol-d6	6.53	99	1392045	92.38	ng	0.00
23) Nitrobenzene-d5	7.45	82	915024	57.85	ng	-0.01
42) 2,4,6-Tribromophenol	10.72	330	503679	84.21	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1771711	56.18	ng	0.00
79) Terphenyl-d14	13.00	244	1815336	59.54	ng	0.00
Target Compounds						
10) Phenol	6.54	94	61134	3.733	ng	Qvalue # 72
50) Dimethylphthalate	9.63	163	185674	5.375	ng	99
71) Phenanthrene	11.43	178	149972	3.243	ng	98
75) Fluoranthene	12.62	202	262791	5.535	ng	98
78) Pyrene	12.85	202	210541	5.287	ng	97
81) Benzo(a)anthracene	14.04	228	99025	2.973	ng	100
83) Chrysene	14.07	228	92559	2.914	ng	94
86) Indeno(1,2,3-cd)pyrene	17.03	276	48622	2.100	ng	94
88) Benzo(b)fluoranthene	15.10	252	98935m	3.534	ng	
90) Benzo(a)pyrene	15.47	252	62911	2.556	ng	97
92) Benzo(g,h,i)perylene	17.47	276	43748	2.259	ng	96

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

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