

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
 Data File : BF112758.D
 Acq On : 28 Feb 2019 13:21
 Operator : JU/SJ
 Sample : K1596-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-15

Manual Integrations
 APPROVED

Sohil
 3/1/2019 2:59:55 PM

Quant Time: Mar 01 05:54:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 27 16:08:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	163464	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	557556	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	239334	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	494733	20.00	ng	0.00
76) Chrysene-d12	13.88	240	433432	20.00	ng	0.00
87) Perylene-d12	15.29	264	296261	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	964479	91.78	ng	0.00
7) Phenol-d6	6.37	99	1161949	88.02	ng	0.00
23) Nitrobenzene-d5	7.30	82	682441	73.24	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	289313	113.87	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	972830	66.96	ng	0.00
79) Terphenyl-d14	12.83	244	1108736	49.59	ng	0.00
Target Compounds						
10) Phenol	6.39	94	43337	2.813	ng	92
50) Dimethylphthalate	9.49	163	76574	4.339	ng	99
71) Phenanthrene	11.27	178	134224	5.453	ng	98
75) Fluoranthene	12.46	202	295625	11.210	ng	97
78) Pyrene	12.69	202	295135	8.077	ng	99
81) Benzo(a)anthracene	13.87	228	150358	5.005	ng	96
83) Chrysene	13.90	228	147659m	5.018	ng	
86) Indeno(1,2,3-cd)pyrene	16.66	276	53503	2.133	ng	97
88) Benzo(b)fluoranthene	14.89	252	143985m	7.514	ng	
90) Benzo(a)pyrene	15.23	252	86743	5.118	ng	# 89
92) Benzo(a,h,i)perylene	17.06	276	56417	3.885	ng	# 91

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

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