

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011819\
 Data File : BF112116.D
 Acq On : 18 Jan 2019 10:09
 Operator : JU/SJ
 Sample : K1157-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-02-011519

Manual Integrations
 APPROVED

Sohil
 1/21/2019 12:00:34 PM

Quant Time: Jan 18 12:20:19 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 16 18:22:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	299460	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1118414	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	525382	20.00	ng	-0.01
64) Phenanthrene-d10	11.38	188	887820	20.00	ng	0.00
76) Chrysene-d12	14.02	240	723121	20.00	ng	-0.01
87) Perylene-d12	15.50	264	789062	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	1587605	96.23	ng	0.00
7) Phenol-d6	6.50	99	1978479	95.84	ng	0.00
23) Nitrobenzene-d5	7.42	82	1238016	76.48	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	534650	94.33	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	2285787	63.79	ng	0.00
79) Terphenyl-d14	12.96	244	1967417	57.88	ng	-0.01
Target Compounds						
10) Phenol	6.52	94	71284	3.160	ng	Qvalue # 74
50) Dimethylphthalate	9.60	163	224669	5.864	ng	98
71) Phenanthrene	11.40	178	597472	13.372	ng	97
75) Fluoranthene	12.59	202	762788	16.169	ng	99
78) Pyrene	12.82	202	770830	16.157	ng	97
81) Benzo(a)anthracene	14.01	228	350327	8.449	ng	96
83) Chrysene	14.04	228	346938	8.843	ng	97
86) Indeno(1,2,3-cd)pyrene	16.97	276	191864	5.301	ng	96
88) Benzo(b)fluoranthene	15.06	252	421440m	9.409	ng	
89) Benzo(k)fluoranthene	15.09	252	146830m	3.387	ng	
90) Benzo(a)pyrene	15.43	252	315246	7.675	ng	# 95
92) Benzo(a,h,i)perylene	17.42	276	187113	5.206	ng	99

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

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