

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011819\
 Data File : BF112123.D
 Acq On : 18 Jan 2019 13:19
 Operator : JU/SJ
 Sample : K1157-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-06-011519

Manual Integrations
 APPROVED

Sohil
 1/21/2019 12:00:43 PM

Quant Time: Jan 18 14:32:56 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	266282	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	951720	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	417343	20.00	ng	-0.01
64) Phenanthrene-d10	11.38	188	661747	20.00	ng	0.00
76) Chrysene-d12	14.02	240	591155	20.00	ng	-0.01
87) Perylene-d12	15.49	264	640150	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1906612	129.96	ng	0.00
7) Phenol-d6	6.50	99	2347798	127.91	ng	0.00
23) Nitrobenzene-d5	7.42	82	1494385	108.48	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	577862	128.35	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	2676715	94.03	ng	0.00
79) Terphenyl-d14	12.96	244	2164269	77.89	ng	-0.01
Target Compounds						
10) Phenol	6.52	94	47243	2.355	ng	82
50) Dimethylphthalate	9.60	163	258132	8.481	ng	97
71) Phenanthrene	11.40	178	113594	3.411	ng	95
75) Fluoranthene	12.59	202	191629	5.450	ng	100
78) Pyrene	12.82	202	181821	4.662	ng	98
81) Benzo(a)anthracene	14.01	228	100280	2.959	ng	97
83) Chrysene	14.04	228	108106m	3.371	ng	
88) Benzo(b)fluoranthene	15.06	252	141426m	3.892	ng	
90) Benzo(a)pyrene	15.43	252	95385	2.862	ng	96

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

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