

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010319\
 Data File : BF111840.D
 Acq On : 4 Jan 2019 00:00
 Operator : JU/SJ
 Sample : K1019-01
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEST-PIT-D-11-12-A

Manual Integrations
 APPROVED

Sohil
 1/4/2019 11:20:50 AM

Quant Time: Jan 04 09:56:25 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	242329	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	894050	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	406945	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	549710	20.00	ng	0.01
76) Chrysene-d12	14.07	240	477272	20.00	ng	0.02
87) Perylene-d12	15.56	264	253343	20.00	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	1390034	97.77	ng	0.02
7) Phenol-d6	6.55	99	1822650	100.74	ng	0.02
23) Nitrobenzene-d5	7.46	82	1107906	72.13	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	334004	69.46	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	2095687	75.06	ng	0.00
79) Terphenyl-d14	13.01	244	1246638	49.53	ng	0.01
Target Compounds						
10) Phenol	6.56	94	86991	4.261	ng	98
50) Dimethylphthalate	9.65	163	316744	10.644	ng	98
71) Phenanthrene	11.45	178	154562	5.302	ng	93
75) Fluoranthene	12.64	202	193512	6.555	ng	94
78) Pyrene	12.87	202	160173	4.752	ng	95
81) Benzo(a)anthracene	14.06	228	92081	3.381	ng	99
83) Chrysene	14.09	228	80832m	3.019	ng	
88) Benzo(b)fluoranthene	15.12	252	48774m	3.294	ng	
90) Benzo(a)pyrene	15.49	252	36309	2.699	ng	98

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

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