

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010419\
 Data File : BF111855.D
 Acq On : 4 Jan 2019 20:03
 Operator : JU/SJ
 Sample : K1041-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OS-1-5-COMP

Manual Integrations
 APPROVED

Sohil
 1/7/2019 11:03:59 AM

Quant Time: Jan 05 00:15:27 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	6.89	152	159383	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	582188	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	267822	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	403358	20.00	ng	0.00
76) Chrysene-d12	14.06	240	354676	20.00	ng	0.00
87) Perylene-d12	15.54	264	318252	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	926759	99.11	ng	0.02
7) Phenol-d6	6.54	99	1215292	102.13	ng	0.02
23) Nitrobenzene-d5	7.45	82	775069	77.49	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	309254	97.73	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1386022	75.43	ng	0.00
79) Terphenyl-d14	13.00	244	1238970	66.25	ng	0.00
Target Compounds						
10) Phenol	6.55	94	61586	4.587	ng	86
50) Dimethylphthalate	9.64	163	131003	6.689	ng	98
71) Phenanthrene	11.44	178	63213	2.955	ng	93
75) Fluoranthene	12.63	202	142454	6.576	ng	95
78) Pyrene	12.86	202	126007	5.031	ng	95
81) Benzo(a)anthracene	14.05	228	59558	2.942	ng	98
83) Chrysene	14.08	228	51862	2.607	ng	97
88) Benzo(b)fluoranthene	15.10	252	64721m	3.480	ng	
90) Benzo(a)pyrene	15.47	252	44862	2.655	ng	98

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

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