

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010819\
 Data File : BF111925.D
 Acq On : 8 Jan 2019 23:28
 Operator : JU/SJ
 Sample : K1075-19 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BASIN-10

Manual Integrations
 APPROVED

Sohil
 1/9/2019 10:36:51 AM

Quant Time: Jan 09 01:39:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 07 14:33:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	116577	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	381512	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	194790	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	206797	20.00	ng	0.00
76) Chrysene-d12	14.06	240	140596	20.00	ng	0.00
87) Perylene-d12	15.54	264	209078	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	335699	49.27	ng	0.00
7) Phenol-d6	6.53	99	376206	42.84	ng	0.00
23) Nitrobenzene-d5	7.44	82	242797	34.01	ng	-0.02
42) 2,4,6-Tribromophenol	10.72	330	59197	23.36	ng	0.00
45) 2-Fluorobiphenyl	9.24	172	376983	28.22	ng	-0.01
79) Terphenyl-d14	12.99	244	251441	33.95	ng	0.00
Target Compounds						
10) Phenol	6.54	94	21602	2.263	ng	Qvalue # 75
31) Naphthalene	8.19	128	53692	2.965	ng	99
50) Dimethylphthalate	9.63	163	50693	3.464	ng	98
52) Acenaphthene	9.96	154	95270	7.817	ng	98
55) Dibenzofuran	10.13	168	68443	3.861	ng	98
58) Fluorene	10.47	166	71424	5.538	ng	96
71) Phenanthrene	11.44	178	601984	54.447	ng	100
72) Anthracene	11.49	178	179876	16.026	ng	98
73) Carbazole	11.64	167	63096	5.779	ng	98
75) Fluoranthene	12.63	202	829484	73.080	ng	99
78) Pyrene	12.86	202	644263	66.598	ng	99
80) Butylbenzylphthalate	13.47	149	29009	7.093	ng	96
81) Benzo(a)anthracene	14.04	228	299919	37.066	ng	99
83) Chrysene	14.09	228	232629	30.153	ng	97
84) Bis(2-ethylhexyl)phthalate	14.03	149	58913	10.969	ng	95
86) Indeno(1,2,3-cd)pyrene	17.05	276	162386m	28.871	ng	
88) Benzo(b)fluoranthene	15.11	252	459143m	36.711	ng	
89) Benzo(k)fluoranthene	15.13	252	141984m	12.077	ng	
90) Benzo(a)pyrene	15.48	252	329358	29.950	ng	# 95
91) Dibenzo(a,h)anthracene	17.06	278	44597	5.042	ng	# 85
92) Benzo(g,h,i)perylene	17.50	276	133815	15.468	ng	98

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

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