

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082119\
 Data File : BF116465.D
 Acq On : 21 Aug 2019 15:02
 Operator : HP/JU
 Sample : K4421-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SUNSET-PARK-SB-6-COMP

Manual Integrations
 APPROVED

Jagrut
 8/22/2019 1:57:56 PM

Quant Time: Aug 21 16:58:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 19 06:03:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	95245	20.00	ng	-0.02
21) Naphthalene-d8	8.07	136	373448	20.00	ng	-0.02
39) Acenaphthene-d10	9.83	164	194978	20.00	ng	-0.02
64) Phenanthrene-d10	11.32	188	298439	20.00	ng	-0.02
76) Chrysene-d12	13.97	240	213337	20.00	ng	-0.01
87) Perylene-d12	15.43	264	206202	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	613147	107.77	ng	-0.02
7) Phenol-d6	6.45	99	813963	113.39	ng	-0.02
23) Nitrobenzene-d5	7.36	82	496795	76.79	ng	-0.02
42) 2,4,6-Tribromophenol	10.62	330	164582	87.06	ng	-0.02
45) 2-Fluorobiphenyl	9.15	172	806298	77.46	ng	-0.02
79) Terphenyl-d14	12.90	244	752372	74.31	ng	-0.02

Target Compounds				Qvalue		
31) Naphthalene	8.10	128	67802	3.858	ng	97
38) 1-Methylnaphthalene	8.89	142	28358	2.590	ng	# 95
46) 1,1'-Biphenyl	9.25	154	31019	2.253	ng	98
49) Acenaphthylene	9.69	152	308240	18.441	ng	98
50) Dimethylphthalate	9.55	163	144450	10.881	ng	99
52) Acenaphthene	9.86	154	116451	11.357	ng	97
58) Fluorene	10.38	166	76240	6.645	ng	97
71) Phenanthrene	11.35	178	381027	24.974	ng	99
72) Anthracene	11.39	178	301089	19.500	ng	99
75) Fluoranthene	12.54	202	1155491	72.343	ng	96
78) Pyrene	12.77	202	1751349	108.904	ng	99
81) Benzo(a)anthracene	13.96	228	620983	45.541	ng	99
83) Chrysene	14.00	228	563265	42.549	ng	100
84) Bis(2-ethylhexyl)phthalate	13.92	149	22345	2.528	ng	# 92
86) Indeno(1,2,3-cd)pyrene	16.89	276	188150	16.922	ng	95
88) Benzo(b)fluoranthene	15.01	252	569255m	47.745	ng	
89) Benzo(k)fluoranthene	15.03	252	122749m	10.570	ng	
90) Benzo(a)pyrene	15.37	252	306517	28.759	ng	# 97
91) Dibenzo(a,h)anthracene	16.89	278	42266	4.581	ng	# 83
92) Benzo(a,h,i)perylene	17.33	276	241632	26.669	ng	97

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

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