

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082819\
 Data File : BF116623.D
 Acq On : 28 Aug 2019 19:07
 Operator : HP/JU
 Sample : K4568-12
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-19-COMP

Manual Integrations
 APPROVED

Jagrut
 8/29/2019 7:46:58 PM

Quant Time: Aug 29 02:25:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 28 13:34:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	87499	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	331293	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	179625	20.00	ng	-0.01
64) Phenanthrene-d10	11.37	188	317377	20.00	ng	0.00
76) Chrysene-d12	14.00	240	149868	20.00	ng	0.00
87) Perylene-d12	15.46	264	127751	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	781962	154.36	ng	0.00
7) Phenol-d6	6.49	99	1047582	152.53	ng	0.00
23) Nitrobenzene-d5	7.42	82	652832	112.76	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	260988	152.68	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	1072326	99.35	ng	0.00
79) Terphenyl-d14	12.96	244	1091251	118.42	ng	0.00

Target Compounds					Qvalue	
10) Phenol	6.50	94	25237	3.350	ng	98
31) Naphthalene	8.16	128	95745	6.033	ng	99
37) 2-Methylnaphthalene	8.85	142	32777	3.307	ng	97
38) 1-Methylnaphthalene	8.95	142	27517	2.843	ng	94
50) Dimethylphthalate	9.60	163	144138	11.936	ng	99
52) Acenaphthene	9.93	154	53304	5.042	ng	96
55) Dibenzofuran	10.10	168	66004	4.636	ng	97
58) Fluorene	10.44	166	69263	6.411	ng	99
71) Phenanthrene	11.40	178	835607	49.018	ng	99
72) Anthracene	11.45	178	179013	10.502	ng	98
73) Carbazole	11.60	167	65358	4.505	ng	98
75) Fluoranthene	12.59	202	726211	47.500	ng	99
78) Pvrene	12.82	202	686391	49.807	ng	99
81) Benzo(a)anthracene	14.00	228	175999	17.614	ng	97
83) Chrysene	14.03	228	163422	16.707	ng	98
86) Indeno(1,2,3-cd)pyrene	16.90	276	62900	11.492	ng	# 94
88) Benzo(b)fluoranthene	15.04	252	132532m	16.893	ng	
89) Benzo(k)fluoranthene	15.06	252	51146m	6.859	ng	
90) Benzo(a)pvrene	15.40	252	104205	14.883	ng	98
91) Dibenzo(a,h)anthracene	16.91	278	14876m	2.277	ng	
92) Benzo(q,h,i)perylene	17.34	276	63291	9.317	ng	97

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

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