

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012419\
 Data File : BF112208.D
 Acq On : 24 Jan 2019 14:56
 Operator : JU/SJ
 Sample : K1003-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-02-012319

Manual Integrations
 APPROVED

Sohil
 1/25/2019 10:05:00 AM

Quant Time: Jan 24 15:30:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 19 01:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	215749	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	835177	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	412461	20.00	ng	-0.01
64) Phenanthrene-d10	11.37	188	805123	20.00	ng	-0.01
76) Chrysene-d12	14.02	240	549537	20.00	ng	-0.01
87) Perylene-d12	15.49	264	572446	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	1078133	90.70	ng	0.00
7) Phenol-d6	6.50	99	1359119	91.39	ng	0.00
23) Nitrobenzene-d5	7.41	82	853278	70.59	ng	-0.02
42) 2,4,6-Tribromophenol	10.68	330	435993	97.99	ng	-0.01
45) 2-Fluorobiphenyl	9.21	172	1698224	60.36	ng	0.00
79) Terphenyl-d14	12.96	244	1634819	63.29	ng	-0.01
Target Compounds						Qvalue
10) Phenol	6.51	94	70552	4.341	ng	79
31) Naphthalene	8.15	128	143778	3.785	ng	97
50) Dimethylphthalate	9.60	163	130704	4.345	ng	100
52) Acenaphthene	9.92	154	89513	3.714	ng	96
55) Dibenzofuran	10.09	168	145143	3.982	ng	98
58) Fluorene	10.43	166	189073	6.942	ng	97
71) Phenanthrene	11.40	178	990976	24.457	ng	98
72) Anthracene	11.45	178	321811	7.840	ng	98
73) Carbazole	11.60	167	144589	3.517	ng	99
75) Fluoranthene	12.59	202	752096	17.580	ng	98
78) Pyrene	12.82	202	571904	15.774	ng	98
81) Benzo(a)anthracene	14.00	228	245170	7.781	ng	98
83) Chrysene	14.04	228	216819	7.272	ng	96
86) Indeno(1,2,3-cd)pyrene	16.97	276	118642	4.313	ng	95
88) Benzo(b)fluoranthene	15.06	252	284241m	8.747	ng	
89) Benzo(k)fluoranthene	15.09	252	77272m	2.457	ng	
90) Benzo(a)pyrene	15.43	252	197323	6.622	ng	99
92) Benzo(g,h,i)perylene	17.42	276	117936	4.523	ng	97

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

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