

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011819\
 Data File : BF112127.D
 Acq On : 18 Jan 2019 16:32
 Operator : JU/SJ
 Sample : K1156-10DL 2X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-7DL

Manual Integrations
 APPROVED

Sohil
 1/21/2019 12:00:49 PM

Quant Time: Jan 18 23:50:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 18 23:33:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	285028	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1112644	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	530949	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	920414	20.00	ng	0.00
76) Chrysene-d12	14.02	240	670571	20.00	ng	0.00
87) Perylene-d12	15.50	264	720607	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	713686	45.45	ng	0.00
7) Phenol-d6	6.50	99	917916	46.72	ng	0.00
23) Nitrobenzene-d5	7.41	82	550874	34.21	ng	-0.02
42) 2,4,6-Tribromophenol	10.68	330	250906	43.81	ng	-0.01
45) 2-Fluorobiphenyl	9.21	172	1049897	28.99	ng	0.00
79) Terphenyl-d14	12.96	244	883797	28.04	ng	0.00
Target Compounds						
10) Phenol	6.51	94	51380	2.393	ng	Qvalue # 75
31) Naphthalene	8.16	128	135097	2.670	ng	98
37) 2-Methylnaphthalene	8.85	142	77478	2.255	ng	97
38) 1-Methylnaphthalene	8.95	142	69661	2.109	ng	99
50) Dimethylphthalate	9.60	163	94331	2.436	ng	# 99
52) Acenaphthene	9.92	154	182087	5.868	ng	98
55) Dibenzofuran	10.10	168	211728	4.512	ng	98
58) Fluorene	10.44	166	242646	6.921	ng	99
71) Phenanthrene	11.41	178	2385256	51.495	ng	99
72) Anthracene	11.46	178	773443	16.483	ng	98
73) Carbazole	11.61	167	189331	4.028	ng	99
75) Fluoranthene	12.60	202	2239027	45.782	ng	99
78) Pvrene	12.83	202	1899976	42.946	ng	100
81) Benzo(a)anthracene	14.01	228	945605	24.594	ng	99
83) Chrysene	14.05	228	771886	21.217	ng	97
86) Indeno(1,2,3-cd)pyrene	16.97	276	362685	10.806	ng	95
88) Benzo(b)fluoranthene	15.07	252	973234m	23.793	ng	
89) Benzo(k)fluoranthene	15.09	252	255375m	6.450	ng	
90) Benzo(a)pvrene	15.43	252	691696	18.439	ng	98
91) Dibenzo(a,h)anthracene	16.98	278	100561m	3.018	ng	
92) Benzo(g,h,i)perylene	17.42	276	313748	9.559	ng	98

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

