

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032919\
 Data File : BF113454.D
 Acq On : 30 Mar 2019 11:35
 Operator : JU/SJ
 Sample : K2153-15
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CONC-8

Manual Integrations
 APPROVED

Sohil
 4/1/2019 5:46:48 PM

Quant Time: Apr 01 05:04:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	148691	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	528852	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	249260	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	466602	20.00	ng	0.00
76) Chrysene-d12	13.86	240	326682	20.00	ng	0.02
87) Perylene-d12	15.27	264	289141	20.00	ng	0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	1340884	147.45	ng	0.01
7) Phenol-d6	6.36	99	1746452	151.81	ng	0.00
23) Nitrobenzene-d5	7.27	82	1061462	119.13	ng	0.00
42) 2,4,6-Tribromophenol	10.53	330	464994	151.02	ng	0.00
45) 2-Fluorobiphenyl	9.07	172	1888139	127.56	ng	0.00
79) Terphenyl-d14	12.81	244	1909345	128.83	ng	0.01

Target Compounds				Qvalue		
10) Phenol	6.37	94	42341	3.224	ng	# 70
31) Naphthalene	8.01	128	88865	3.439	ng	99
37) 2-Methylnaphthalene	8.70	142	53242	3.302	ng	98
38) 1-Methylnaphthalene	8.80	142	71518	4.611	ng	97
49) Acenaphthylene	9.60	152	85886	3.314	ng	# 97
50) Dimethylphthalate	9.46	163	136202	7.373	ng	100
52) Acenaphthene	9.77	154	225267	15.434	ng	99
55) Dibenzofuran	9.94	168	142302	6.485	ng	# 96
58) Fluorene	10.28	166	275739	16.177	ng	99
71) Phenanthrene	11.26	178	3018410	130.264	ng	98
72) Anthracene	11.30	178	928533	39.451	ng	100
73) Carbazole	11.45	167	236159	10.515	ng	99
75) Fluoranthene	12.45	202	4044635	154.405	ng	97
78) Pyrene	12.68	202	4056535	179.192	ng	95
81) Benzo(a)anthracene	13.85	228	2218261	118.625	ng	96
83) Chrysene	13.89	228	2017218	106.204	ng	95
84) Bis(2-ethylhexyl)phthalate	13.84	149	572949	46.846	ng	# 98
86) Indeno(1,2,3-cd)pyrene	16.64	276	980317	60.139	ng	95
88) Benzo(b)fluoranthene	14.88	252	2062274m	114.817	ng	
89) Benzo(k)fluoranthene	14.90	252	728813m	46.205	ng	
90) Benzo(a)pyrene	15.22	252	1546154	97.099	ng	97
91) Dibenz(a,h)anthracene	16.64	278	247333m	17.964	ng	
92) Benzo(g,h,i)perylene	17.06	276	983289	70.830	ng	97

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

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