

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120718\
 Data File : BF111210.D
 Acq On : 8 Dec 2018 2:29
 Operator : JU/SJ
 Sample : J6260-09
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-03-COMP

Manual Integrations
 APPROVED

Sohil
 12/11/2018 8:44:55 AM

Quant Time: Dec 08 03:40:06 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 05 17:30:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	417299	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	1462661	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	517325	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	916106	20.00	ng	0.00
76) Chrysene-d12	13.96	240	672280	20.00	ng	0.00
87) Perylene-d12	15.40	264	505980	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	2327301	87.03	ng	0.00
7) Phenol-d6	6.43	99	3006627	87.28	ng	0.00
23) Nitrobenzene-d5	7.36	82	1951990	79.01	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	420883	102.35	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	2761280	101.61	ng	0.00
79) Terphenyl-d14	12.91	244	1595115	48.59	ng	0.00

Target Compounds

						Qvalue
10) Phenol	6.45	94	190905	5.082	ng	96
50) Dimethylphthalate	9.55	163	171009	4.750	ng	98
52) Acenaphthene	9.87	154	82436	2.713	ng	96
58) Fluorene	10.39	166	90010	3.050	ng	99
71) Phenanthrene	11.34	178	1454032	32.997	ng	96
72) Anthracene	11.40	178	341925	7.698	ng	94
73) Carbazole	11.54	167	157535	3.408	ng	94
75) Fluoranthene	12.53	202	2178008	50.510	ng	99
78) Pyrene	12.76	202	1815545	34.664	ng	97
81) Benzo(a)anthracene	13.95	228	965564	24.514	ng	99
83) Chrysene	13.99	228	947578	24.233	ng	97
84) Bis(2-ethylhexyl)phthalate	13.95	149	174055	5.695	ng	99
86) Indeno(1,2,3-cd)pyrene	16.80	276	314283	10.010	ng	# 93
88) Benzo(b)fluoranthene	14.98	252	906823m	31.629	ng	
89) Benzo(k)fluoranthene	15.01	252	326792m	11.790	ng	
90) Benzo(a)pyrene	15.33	252	613766	23.339	ng	98
91) Dibenzo(a,h)anthracene	16.81	278	80569m	3.651	ng	
92) Benzo(g,h,i)perylene	17.23	276	256297	11.557	ng	99

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

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