

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122818\
 Data File : BF111790.D
 Acq On : 28 Dec 2018 15:57
 Operator : JU/SJ
 Sample : J6195-11 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-01-122618-A

Manual Integrations
 APPROVED

Sohil
 1/2/2019 2:51:33 PM

Quant Time: Dec 30 21:25:07 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	165535	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	625092	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	282279	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	441701	20.00	ng	0.01
76) Chrysene-d12	14.07	240	366414	20.00	ng	0.02
87) Perylene-d12	15.56	264	295968	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	452458	46.59	ng	0.02
7) Phenol-d6	6.54	99	592623	47.95	ng	0.02
23) Nitrobenzene-d5	7.46	82	372749	34.71	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	141331	42.37	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	679440	35.08	ng	0.00
79) Terphenyl-d14	13.01	244	527177	27.28	ng	0.01
Target Compounds						
10) Phenol	6.55	94	40416	2.898	ng	97
31) Naphthalene	8.20	128	141724	4.719	ng	96
37) 2-Methylnaphthalene	8.90	142	68722	3.517	ng	100
38) 1-Methylnaphthalene	9.00	142	47777m	2.546	ng	
50) Dimethylphthalate	9.65	163	116044	5.622	ng	# 97
52) Acenaphthene	9.97	154	109137	6.419	ng	97
55) Dibenzofuran	10.15	168	134659	5.243	ng	96
58) Fluorene	10.49	166	164911	8.910	ng	99
71) Phenanthrene	11.45	178	1029864	43.964	ng	94
72) Anthracene	11.50	178	346054	14.718	ng	94
73) Carbazole	11.65	167	74718	3.273	ng	93
75) Fluoranthene	12.64	202	1154610	48.674	ng	96
78) Pvrene	12.87	202	959010	37.063	ng	97
81) Benzo(a)anthracene	14.06	228	481492	23.026	ng	99
83) Chrysene	14.09	228	398044	19.367	ng	97
86) Indeno(1,2,3-cd)pyrene	17.06	276	164939	9.440	ng	# 84
88) Benzo(b)fluoranthene	15.12	252	448579m	25.933	ng	
89) Benzo(k)fluoranthene	15.14	252	162160m	9.847	ng	
90) Benzo(a)pvrene	15.50	252	319902m	20.357	ng	
91) Dibenzo(a,h)anthracene	17.07	278	38308	2.784	ng	# 90
92) Benzo(g,h,i)perylene	17.52	276	144412	10.747	ng	# 90

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

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