

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120518\
 Data File : BF111114.D
 Acq On : 5 Dec 2018 18:31
 Operator : JU/SJ
 Sample : J5775-03 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-113018

Manual Integrations
 APPROVED

mohammad
 12/6/2018 12:11:48 PM

Quant Time: Dec 05 18:59:50 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	551054	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	2201872	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	994905	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	1486154	20.00	ng	0.00
76) Chrysene-d12	13.96	240	982938	20.00	ng	0.00
87) Perylene-d12	15.39	264	777360	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	1610640	45.61	ng	-0.01
7) Phenol-d6	6.42	99	2050544	45.08	ng	-0.01
23) Nitrobenzene-d5	7.36	82	1229333	33.05	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	321038	40.59	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	2119257	30.15	ng	0.00
79) Terphenyl-d14	12.92	244	1466013	30.54	ng	0.01
Target Compounds						
10) Phenol	6.43	94	124439	2.509	ng	92
50) Dimethylphthalate	9.55	163	147588	2.131	ng	# 96
52) Acenaphthene	9.87	154	143817	2.461	ng	97
58) Fluorene	10.38	166	225234	3.968	ng	98
71) Phenanthrene	11.35	178	1674387	23.422	ng	96
72) Anthracene	11.39	178	491333	6.819	ng	95
73) Carbazole	11.54	167	180779	2.411	ng	95
75) Fluoranthene	12.54	202	2104749	30.089	ng	83
78) Pyrene	12.77	202	3287921	42.936	ng	99
81) Benzo(a)anthracene	13.95	228	1989050	34.538	ng	99
83) Chrysene	13.99	228	1903648	33.297	ng	98
84) Bis(2-ethylhexyl)phthalate	13.95	149	478870	10.717	ng	100
86) Indeno(1,2,3-cd)pyrene	16.80	276	984847	21.455	ng	92
88) Benzo(b)fluoranthene	14.98	252	2382804m	54.095	ng	
89) Benzo(k)fluoranthene	15.00	252	775048m	18.201	ng	
90) Benzo(a)pyrene	15.33	252	1597780	39.546	ng	98
91) Dibenzo(a,h)anthracene	16.82	278	241628m	7.127	ng	
92) Benzo(g,h,i)perylene	17.23	276	919499	26.987	ng	99

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

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