

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120518\
 Data File : BM017913.D
 Acq On : 04 Dec 2018 21:35
 Operator : JU/SJ
 Sample : J5771-07 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EO-02-112918-A

Manual Integrations
 APPROVED

mohammad
 12/5/2018 4:50:06 PM

Quant Time: Dec 05 01:09:56 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM111218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 12 18:37:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	272442	20.00	ng	-0.04
21) Naphthalene-d8	10.35	136	1054546	20.00	ng	-0.04
39) Acenaphthene-d10	14.23	164	526200	20.00	ng	-0.04
64) Phenanthrene-d10	16.99	188	936949	20.00	ng	-0.03
76) Chrysene-d12	21.19	240	890655	20.00	ng	-0.02
87) Perylene-d12	23.38	264	977390	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.21	112	814365	50.45	ng	-0.03
7) Phenol-d6	6.77	99	1104977	49.01	ng	-0.03
23) Nitrobenzene-d5	8.73	82	648415	31.74	ng	-0.04
42) 2,4,6-Tribromophenol	15.73	330	320980	42.45	ng	-0.03
45) 2-Fluorobiphenyl	12.84	172	1341664	33.08	ng	-0.04
79) Terphenyl-d14	19.63	244	1255100	31.94	ng	-0.02
Target Compounds						
10) Phenol	6.80	94	63705	2.692	ng	Qvalue 96
50) Dimethylphthalate	13.69	163	125028	2.922	ng	98
58) Fluorene	15.29	166	120198	2.972	ng	97
71) Phenanthrene	17.03	178	903164	17.764	ng	99
72) Anthracene	17.12	178	302123	6.107	ng	100
75) Fluoranthene	19.06	202	1172957	20.390	ng	98
78) Pyrene	19.42	202	1056862	19.232	ng	98
81) Benzo(a)anthracene	21.18	228	515023	10.037	ng	98
83) Chrysene	21.23	228	432090	8.670	ng	97
86) Indeno(1,2,3-cd)pyrene	25.55	276	262003	6.579	ng	97
88) Benzo(b)fluoranthene	22.73	252	556894m	9.729	ng	
89) Benzo(k)fluoranthene	22.77	252	197615m	3.416	ng	
90) Benzo(a)pyrene	23.28	252	426016	8.163	ng	# 93
92) Benzo(a,h,i)perylene	26.21	276	244924	5.936	ng	# 93

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

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