

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120518\
 Data File : BM017912.D
 Acq On : 04 Dec 2018 20:59
 Operator : JU/SJ
 Sample : J5770-15 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OR-03-112918-B

Manual Integrations
 APPROVED

mohammad
 12/5/2018 4:50:04 PM

Quant Time: Dec 05 01:07:53 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	290455	20.00	ng	-0.04
21) Naphthalene-d8	10.35	136	1120798	20.00	ng	-0.04
39) Acenaphthene-d10	14.23	164	554831	20.00	ng	-0.04
64) Phenanthrene-d10	16.99	188	1026992	20.00	ng	-0.03
76) Chrysene-d12	21.19	240	946937	20.00	ng	-0.02
87) Perylene-d12	23.38	264	1040240	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.21	112	873558	50.76	ng	-0.03
7) Phenol-d6	6.77	99	1174177	48.85	ng	-0.03
23) Nitrobenzene-d5	8.73	82	691232	31.84	ng	-0.04
42) 2,4,6-Tribromophenol	15.73	330	348725	43.74	ng	-0.03
45) 2-Fluorobiphenyl	12.84	172	1522675	35.60	ng	-0.04
79) Terphenyl-d14	19.63	244	1484822	35.54	ng	-0.02
Target Compounds						
10) Phenol	6.80	94	69225	2.744	ng	Qvalue 94
50) Dimethylphthalate	13.70	163	106304	2.357	ng	99
71) Phenanthrene	17.03	178	254141	4.560	ng	99
75) Fluoranthene	19.05	202	461071	7.312	ng	99
78) Pyrene	19.42	202	459639	7.867	ng	99
81) Benzo(a)anthracene	21.17	228	270960	4.967	ng	95
83) Chrysene	21.23	228	265745m	5.015	ng	
86) Indeno(1,2,3-cd)pyrene	25.54	276	207398	4.898	ng	96
88) Benzo(b)fluoranthene	22.73	252	402948m	6.614	ng	
89) Benzo(k)fluoranthene	22.76	252	126883m	2.061	ng	
90) Benzo(a)pyrene	23.28	252	279626	5.034	ng	# 94
92) Benzo(a,h,i)perylene	26.21	276	199026	4.532	ng	97

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

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