

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122018\
 Data File : BM018340.D
 Acq On : 20 Dec 2018 21:10
 Operator : JU/SJ
 Sample : J6386-10
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS006-0612-01

Manual Integrations
 APPROVED

Sohil
 12/21/2018 2:49:42 PM

Quant Time: Dec 21 02:52:01 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 15 21:39:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.49	152	156662	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	557254	20.00	ng	-0.02
39) Acenaphthene-d10	14.14	164	273236	20.00	ng	-0.02
64) Phenanthrene-d10	16.91	188	552142	20.00	ng	-0.01
76) Chrysene-d12	21.14	240	638387	20.00	ng	0.00
87) Perylene-d12	23.30	264	692468	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	1072592	114.37	ng	0.00
7) Phenol-d6	6.70	99	1365871	104.78	ng	0.00
23) Nitrobenzene-d5	8.65	82	889285	81.86	ng	-0.02
42) 2,4,6-Tribromophenol	15.66	330	397434	99.10	ng	-0.01
45) 2-Fluorobiphenyl	12.75	172	1556369	73.78	ng	-0.02
79) Terphenyl-d14	19.57	244	1854399	65.70	ng	0.00
Target Compounds						
10) Phenol	6.72	94	94004	6.810	ng	91
49) Acenaphthylene	13.86	152	78345	2.873	ng	# 93
50) Dimethylphthalate	13.62	163	128889	5.674	ng	99
71) Phenanthrene	16.95	178	168782	5.628	ng	97
75) Fluoranthene	18.99	202	128636	3.689	ng	95
78) Pyrene	19.35	202	230316	6.055	ng	100
81) Benzo(a)anthracene	21.12	228	97340	2.610	ng	# 84
83) Chrysene	21.17	228	99546	2.775	ng	# 93
88) Benzo(b)fluoranthene	22.67	252	113066m	2.907	ng	
90) Benzo(a)pyrene	23.21	252	103368	2.794	ng	# 79
92) Benzo(g,h,i)perylene	26.09	276	99494	3.071	ng	# 92

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

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