

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122118\
 Data File : BM018364.D
 Acq On : 21 Dec 2018 19:27
 Operator : JU/SJ
 Sample : J6386-19
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS008-0006-01

Manual Integrations
 APPROVED

Sohil
 12/26/2018 3:20:06 PM

Quant Time: Dec 22 05:01:20 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	7.48	152	117490	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	426306	20.00	ng	-0.02
39) Acenaphthene-d10	14.14	164	209874	20.00	ng	-0.02
64) Phenanthrene-d10	16.91	188	399158	20.00	ng	-0.01
76) Chrysene-d12	21.14	240	429872	20.00	ng	0.00
87) Perylene-d12	23.31	264	486522	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	812389	115.51	ng	0.00
7) Phenol-d6	6.70	99	1045798	106.98	ng	0.00
23) Nitrobenzene-d5	8.65	82	710277	85.46	ng	-0.02
42) 2,4,6-Tribromophenol	15.66	330	280322	91.00	ng	-0.01
45) 2-Fluorobiphenyl	12.75	172	1076044	66.41	ng	-0.02
79) Terphenyl-d14	19.57	244	1058577	55.69	ng	0.00
Target Compounds						
10) Phenol	6.73	94	85667	8.275	ng	88
20) 3+4-Methylphenols	8.29	107	19719	2.036	ng	92
49) Acenaphthylene	13.86	152	85788	4.096	ng	97
50) Dimethylphthalate	13.62	163	84287	4.831	ng	97
71) Phenanthrene	16.96	178	228568	10.542	ng	98
72) Anthracene	17.04	178	62016	2.883	ng	97
75) Fluoranthene	18.99	202	162545	6.448	ng	97
78) Pyrene	19.36	202	292417	11.416	ng	99
81) Benzo(a)anthracene	21.13	228	114512m	4.560	ng	
83) Chrysene	21.18	228	125326	5.189	ng	96
86) Indeno(1,2,3-cd)pyrene	25.45	276	76649	3.186	ng	97
88) Benzo(b)fluoranthene	22.67	252	122636m	4.487	ng	
90) Benzo(a)pyrene	23.21	252	133918	5.152	ng	# 90
92) Benzo(a,h,i)perylene	26.11	276	111732	4.909	ng	# 93

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

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