

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122018\
 Data File : BM018336.D
 Acq On : 20 Dec 2018 18:45
 Operator : JU/SJ
 Sample : J6357-08
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-09A

Manual Integrations
 APPROVED

Sohil
 12/21/2018 2:49:41 PM

Quant Time: Dec 21 01:57:20 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	143471	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	587640	20.00	ng	-0.02
39) Acenaphthene-d10	14.14	164	332189	20.00	ng	-0.02
64) Phenanthrene-d10	16.90	188	690564	20.00	ng	-0.02
76) Chrysene-d12	21.13	240	572434	20.00	ng	-0.01
87) Perylene-d12	23.29	264	578768	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	1121926	130.63	ng	0.00
7) Phenol-d6	6.69	99	1601346	134.14	ng	-0.01
23) Nitrobenzene-d5	8.64	82	1179285	102.94	ng	-0.02
42) 2,4,6-Tribromophenol	15.65	330	466657	95.71	ng	-0.02
45) 2-Fluorobiphenyl	12.74	172	2166877	84.49	ng	-0.02
79) Terphenyl-d14	19.56	244	2396861	94.70	ng	-0.01
Target Compounds						
10) Phenol	6.72	94	124245	9.828	ng	99
50) Dimethylphthalate	13.62	163	172449	6.244	ng	99
71) Phenanthrene	16.94	178	94589	2.522	ng	98
75) Fluoranthene	18.98	202	166062	3.808	ng	100
78) Pyrene	19.35	202	174577	5.118	ng	99
81) Benzo(a)anthracene	21.12	228	114675	3.429	ng	97
83) Chrysene	21.17	228	106522m	3.312	ng	
86) Indeno(1,2,3-cd)pyrene	25.42	276	79194	2.472	ng	95
88) Benzo(b)fluoranthene	22.66	252	153684	4.727	ng	# 95
90) Benzo(a)pyrene	23.20	252	95843	3.100	ng	# 88
92) Benzo(g,h,i)perylene	26.07	276	74052	2.735	ng	96

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

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