

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121918\
 Data File : BM018313.D
 Acq On : 20 Dec 2018 01:13
 Operator : JU/SJ
 Sample : J6379-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS004-0006-01

Manual Integrations
 APPROVED

Sohil
 12/20/2018 2:55:13 PM

Quant Time: Dec 20 03:00:36 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	110771	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	396215	20.00	ng	-0.02
39) Acenaphthene-d10	14.14	164	195482	20.00	ng	-0.02
64) Phenanthrene-d10	16.91	188	402598	20.00	ng	-0.01
76) Chrysene-d12	21.13	240	486459	20.00	ng	-0.02
87) Perylene-d12	23.29	264	542169	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	834281	125.81	ng	0.00
7) Phenol-d6	6.69	99	1099891	119.34	ng	-0.02
23) Nitrobenzene-d5	8.64	82	706511	91.46	ng	-0.02
42) 2,4,6-Tribromophenol	15.66	330	334685	116.65	ng	-0.01
45) 2-Fluorobiphenyl	12.75	172	1221700	80.95	ng	-0.02
79) Terphenyl-d14	19.56	244	1404282	65.29	ng	-0.01
Target Compounds						
10) Phenol	6.72	94	64252	6.583	ng	94
50) Dimethylphthalate	13.62	163	103959	6.397	ng	100
75) Fluoranthene	18.98	202	116397	4.578	ng	99
78) Pyrene	19.34	202	99590	3.436	ng	97
81) Benzo(a)anthracene	21.11	228	88039	3.098	ng	98
83) Chrysene	21.16	228	105914	3.875	ng	96
86) Indeno(1,2,3-cd)pyrene	25.41	276	120697	4.433	ng	95
88) Benzo(b)fluoranthene	22.64	252	207100	6.800	ng	# 93
89) Benzo(k)fluoranthene	22.69	252	73066m	2.409	ng	
90) Benzo(a)pyrene	23.19	252	125485	4.332	ng	99
92) Benzo(g,h,i)perylene	26.06	276	126563	4.990	ng	96

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

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