

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121918\
 Data File : BM018312.D
 Acq On : 20 Dec 2018 00:37
 Operator : JU/SJ
 Sample : J6379-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS003-0006-01

Quant Time: Dec 20 02:58:28 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	107647	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	386431	20.00	ng	-0.02
39) Acenaphthene-d10	14.14	164	190204	20.00	ng	-0.02
64) Phenanthrene-d10	16.90	188	387766	20.00	ng	-0.02
76) Chrysene-d12	21.13	240	448933	20.00	ng	-0.02
87) Perylene-d12	23.28	264	502715	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	709009	110.03	ng	0.00
7) Phenol-d6	6.69	99	916906	102.37	ng	-0.02
23) Nitrobenzene-d5	8.64	82	593349	78.76	ng	-0.02
42) 2,4,6-Tribromophenol	15.65	330	269381	96.49	ng	-0.02
45) 2-Fluorobiphenyl	12.75	172	957180	65.18	ng	-0.02
79) Terphenyl-d14	19.56	244	960398	48.38	ng	-0.01
Target Compounds						
10) Phenol	6.71	94	52259	5.509	ng	81
50) Dimethylphthalate	13.62	163	71446	4.518	ng	99
81) Benzo(a)anthracene	21.11	228	52524	2.003	ng	94
83) Chrysene	21.16	228	56927	2.257	ng	96
86) Indeno(1,2,3-cd)pyrene	25.41	276	86920	3.459	ng	95
88) Benzo(b)fluoranthene	22.64	252	142110	5.032	ng	# 97
90) Benzo(a)pyrene	23.19	252	86727	3.229	ng	# 95
92) Benzo(g,h,i)perylene	26.06	276	92290	3.924	ng	96

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

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