

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122118\
 Data File : BM018370.D
 Acq On : 21 Dec 2018 23:05
 Operator : JU/SJ
 Sample : J6499-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 72-11989

Manual Integrations
 APPROVED

Sohil
 12/26/2018 3:20:14 PM

Quant Time: Dec 22 05:46:06 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	119403	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	483273	20.00	ng	-0.02
39) Acenaphthene-d10	14.14	164	265227	20.00	ng	-0.02
64) Phenanthrene-d10	16.91	188	515149	20.00	ng	-0.01
76) Chrysene-d12	21.14	240	428136	20.00	ng	0.00
87) Perylene-d12	23.30	264	473923	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	879607	123.06	ng	0.00
7) Phenol-d6	6.70	99	1196045	120.39	ng	0.00
23) Nitrobenzene-d5	8.65	82	790739	83.93	ng	-0.02
42) 2,4,6-Tribromophenol	15.66	330	354334	91.02	ng	0.00
45) 2-Fluorobiphenyl	12.75	172	1397072	68.23	ng	-0.02
79) Terphenyl-d14	19.57	244	1445604	76.36	ng	0.00
Target Compounds						
10) Phenol	6.73	94	91556	8.702	ng	84
50) Dimethylphthalate	13.62	163	169341	7.680	ng	99
71) Phenanthrene	16.96	178	71853	2.568	ng	99
75) Fluoranthene	18.99	202	252443	7.760	ng	99
78) Pyrene	19.36	202	224619	8.805	ng	98
81) Benzo(a)anthracene	21.12	228	143947	5.756	ng	98
83) Chrysene	21.17	228	132511	5.509	ng	98
86) Indeno(1,2,3-cd)pyrene	25.44	276	75006	3.130	ng	# 93
88) Benzo(b)fluoranthene	22.66	252	169223	6.357	ng	# 88
89) Benzo(k)fluoranthene	22.70	252	58573m	2.209	ng	
90) Benzo(a)pyrene	23.21	252	123245	4.868	ng	# 93
92) Benzo(a,h,i)perylene	26.10	276	70707	3.189	ng	# 93

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

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