

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080818\
 Data File : BM016253.D
 Acq On : 08 Aug 2018 17:05
 Operator : SJ/JU
 Sample : J4313-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C-10

Manual Integrations
 APPROVED

Sohil
 8/9/2018 2:23:55 PM

Quant Time: Aug 09 04:13:50 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 16:23:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	43553	20.00	ng	0.00
21) Naphthalene-d8	10.95	136	158177	20.00	ng	0.00
38) Acenaphthene-d10	14.77	164	75173	20.00	ng	0.00
63) Phenanthrene-d10	17.51	188	148760	20.00	ng	0.00
75) Chrysene-d12	21.64	240	162236	20.00	ng	0.00
86) Perylene-d12	23.98	264	180191	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	236711	90.28	ng	0.00
7) Phenol-d6	7.30	99	283908	82.92	ng	0.00
23) Nitrobenzene-d5	9.32	82	208361	61.27	ng	0.00
41) 2,4,6-Tribromophenol	16.26	330	74862	78.52	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	375836	60.83	ng	0.00
78) Terphenyl-d14	20.09	244	479429	61.86	ng	0.00
Target Compounds						
49) Dimethylphthalate	14.23	163	28974	4.404	ng	Qvalue # 97
70) Phenanthrene	17.55	178	59409	7.133	ng	97
74) Fluoranthene	19.54	202	97522	10.497	ng	98
77) Pyrene	19.90	202	83914	8.631	ng	98
80) Benzo(a)anthracene	21.62	228	59632	5.968	ng	97
82) Chrysene	21.67	228	50204	5.238	ng	96
85) Indeno(1,2,3-cd)pyrene	26.36	276	37912	3.333	ng	# 84
87) Benzo(b)fluoranthene	23.27	252	75570	7.123	ng	97
88) Benzo(k)fluoranthene	23.32	252	24837m	2.310	ng	
89) Benzo(a)pyrene	23.87	252	50092	4.911	ng	92
91) Benzo(g,h,i)perylene	27.10	276	38688	3.871	ng	# 93

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

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