

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080818\
 Data File : BM016250.D
 Acq On : 08 Aug 2018 15:14
 Operator : SJ/JU
 Sample : J4313-07
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C-8

Manual Integrations
 APPROVED

Sohil
 8/9/2018 2:23:50 PM

Quant Time: Aug 08 17:13:35 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	8.13	152	43399	20.00	ng	0.00
21) Naphthalene-d8	10.95	136	156392	20.00	ng	0.00
38) Acenaphthene-d10	14.77	164	75902	20.00	ng	0.00
63) Phenanthrene-d10	17.50	188	155070	20.00	ng	0.00
75) Chrysene-d12	21.63	240	168074	20.00	ng	0.00
86) Perylene-d12	23.97	264	182930	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	286926	109.82	ng	0.00
7) Phenol-d6	7.30	99	339864	99.61	ng	0.00
23) Nitrobenzene-d5	9.32	82	252894	75.21	ng	0.00
41) 2,4,6-Tribromophenol	16.26	330	98668	102.49	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	468778	75.15	ng	0.00
78) Terphenyl-d14	20.09	244	638629	79.54	ng	0.00
Target Compounds						
49) Dimethylphthalate	14.22	163	35290	5.313	ng	98
70) Phenanthrene	17.54	178	160295	18.463	ng	98
71) Anthracene	17.64	178	34886	4.135	ng	# 94
74) Fluoranthene	19.54	202	280653	28.980	ng	98
77) Pyrene	19.90	202	251464	24.966	ng	99
80) Benzo(a)anthracene	21.62	228	159440	15.402	ng	98
82) Chrysene	21.67	228	135989	13.695	ng	95
85) Indeno(1,2,3-cd)pyrene	26.36	276	98124	8.327	ng	# 95
87) Benzo(b)fluoranthene	23.27	252	197178	18.307	ng	# 95
88) Benzo(k)fluoranthene	23.31	252	66556m	6.097	ng	
89) Benzo(a)pyrene	23.87	252	139083	13.433	ng	97
90) Dibenzo(a,h)anthracene	26.34	278	32006	2.984	ng	# 85
91) Benzo(a,h,i)perylene	27.10	276	96459	9.507	ng	# 91

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

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