

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080818\
 Data File : BM016252.D
 Acq On : 08 Aug 2018 16:28
 Operator : SJ/JU
 Sample : J4283-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TR-04-080318

Manual Integrations
 APPROVED

Sohil
 8/9/2018 4:36:12 PM

Quant Time: Aug 09 13:55:27 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	43214	20.00	ng	0.00
21) Naphthalene-d8	10.95	136	162783	20.00	ng	0.00
38) Acenaphthene-d10	14.77	164	83960	20.00	ng	0.00
63) Phenanthrene-d10	17.50	188	178380	20.00	ng	0.00
75) Chrysene-d12	21.64	240	203799	20.00	ng	0.00
86) Perylene-d12	23.97	264	231042	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	214170	82.33	ng	0.00
7) Phenol-d6	7.30	99	261118	76.86	ng	0.00
23) Nitrobenzene-d5	9.32	82	188811	53.95	ng	0.00
41) 2,4,6-Tribromophenol	16.26	330	80603	75.69	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	336962	48.83	ng	0.00
78) Terphenyl-d14	20.09	244	499878	51.34	ng	0.00
Target Compounds						
49) Dimethylphthalate	14.22	163	35841	4.878	ng	Qvalue # 98
74) Fluoranthene	19.54	202	32800	2.944	ng	97
77) Pyrene	19.90	202	45919	3.760	ng	97
82) Chrysene	21.67	228	25887m	2.150	ng	
87) Benzo(b)fluoranthene	23.27	252	39882	2.932	ng	95
89) Benzo(a)pyrene	23.88	252	31559	2.413	ng	# 86
91) Benzo(g,h,i)perylene	27.11	276	26171m	2.042	ng	

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

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