

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080718\
 Data File : BM016230.D
 Acq On : 07 Aug 2018 23:18
 Operator : SJ/JU
 Sample : J4335-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CINNAMINSON-8-3-18

Manual Integrations
 APPROVED

Sohil
 8/8/2018 5:35:07 PM

Quant Time: Aug 08 02:36:42 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 07 16:15:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	54133	20.00	ng	0.00
21) Naphthalene-d8	10.95	136	227952	20.00	ng	0.00
38) Acenaphthene-d10	14.77	164	141789	20.00	ng	0.00
63) Phenanthrene-d10	17.51	188	329348	20.00	ng	0.00
75) Chrysene-d12	21.64	240	372706	20.00	ng	0.00
86) Perylene-d12	23.98	264	335629	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	380080	116.63	ng	0.00
7) Phenol-d6	7.30	99	515723	121.18	ng	0.00
23) Nitrobenzene-d5	9.32	82	344709	70.33	ng	0.00
41) 2,4,6-Tribromophenol	16.26	330	208854	116.14	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	769031	65.99	ng	0.00
78) Terphenyl-d14	20.09	244	1531428	86.01	ng	0.00
Target Compounds						
49) Dimethylphthalate	14.23	163	66365	5.348	ng	98
74) Fluoranthene	19.54	202	105252	5.117	ng	97
77) Pyrene	19.90	202	111983	5.014	ng	99
80) Benzo(a)anthracene	21.62	228	69242	3.016	ng	97
82) Chrysene	21.67	228	77708m	3.529	ng	
87) Benzo(b)fluoranthene	23.27	252	95831	4.849	ng	97
89) Benzo(a)pyrene	23.88	252	61151	3.219	ng	# 93
91) Benzo(g,h,i)perylene	27.10	276	42953	2.307	ng	95

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

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