

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072220\
 Data File : BM026840.D
 Acq On : 22 Jul 2020 19:53
 Operator : JU/CG
 Sample : L3365-04 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 52-09-57

Manual Integrations
 APPROVED

Sohil
 7/24/2020 12:11:35 PM

Quant Time: Jul 23 02:59:32 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 22 14:16:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.30	152	83433	20.00	ng	0.00
21) Naphthalene-d8	10.05	136	307942	20.00	ng	0.00
39) Acenaphthene-d10	13.96	164	163267	20.00	ng	0.00
64) Phenanthrene-d10	16.72	188	329139	20.00	ng	0.00
76) Chrysene-d12	20.94	240	400637	20.00	ng	0.00
86) Perylene-d12	23.01	264	439382	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.94	112	216434	43.48	ng	0.00
7) Phenol-d6	6.51	99	277332	34.97	ng	0.00
23) Nitrobenzene-d5	8.44	82	195819	27.13	ng	0.00
42) 2,4,6-Tribromophenol	15.47	330	75214	31.63	ng	0.00
45) 2-Fluorobiphenyl	12.57	172	340254	27.71	ng	0.00
79) Terphenyl-d14	19.38	244	472040	23.19	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.44	163	47598	3.408	ng	97
71) Phenanthrene	16.76	178	394617	20.309	ng	99
72) Anthracene	16.85	178	84838	4.386	ng	97
73) Carbazole	17.14	167	37587	2.038	ng	98
75) Fluoranthene	18.79	202	779693	32.579	ng	99
78) Pyrene	19.16	202	692166	27.426	ng	99
81) Benzo(a)anthracene	20.93	228	428032	15.905	ng	99
83) Chrysene	20.98	228	414664	15.831	ng	96
87) Indeno(1,2,3-cd)pyrene	24.99	276	326304	10.432	ng	96
88) Benzo(b)fluoranthene	22.41	252	637877m	22.033	ng	
89) Benzo(k)fluoranthene	22.44	252	165807m	5.839	ng	
90) Benzo(a)pyrene	22.92	252	442408	16.461	ng	96
91) Dibenzo(a,h)anthracene	24.99	278	89055	3.373	ng	# 82
92) Benzo(g,h,i)perylene	25.60	276	328312	13.494	ng	99

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

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