

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072220\
 Data File : BM026841.D
 Acq On : 22 Jul 2020 20:30
 Operator : JU/CG
 Sample : L3360-03RX
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP-2RX

Manual Integrations
 APPROVED

Sohil
 7/24/2020 12:11:36 PM

Quant Time: Jul 23 03:05:26 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	101738	20.00	ng	0.00
21) Naphthalene-d8	10.05	136	355164	20.00	ng	0.00
39) Acenaphthene-d10	13.96	164	185425	20.00	ng	0.00
64) Phenanthrene-d10	16.72	188	392072	20.00	ng	0.00
76) Chrysene-d12	20.94	240	443001	20.00	ng	0.00
86) Perylene-d12	23.01	264	473388	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.95	112	19521	3.22	ng	0.00
7) Phenol-d6	6.51	99	279454	28.90	ng	0.00
23) Nitrobenzene-d5	8.44	82	437057	52.50	ng	0.00
42) 2,4,6-Tribromophenol	15.48	330	5013	1.86	ng	0.00
45) 2-Fluorobiphenyl	12.57	172	735664	52.75	ng	0.00
79) Terphenyl-d14	19.39	244	1064224	47.28	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.44	163	108799	6.860	ng	99
71) Phenanthrene	16.76	178	111265	4.807	ng	98
75) Fluoranthene	18.79	202	155110	5.441	ng	96
78) Pyrene	19.16	202	176938	6.340	ng	99
81) Benzo(a)anthracene	20.93	228	95049	3.194	ng	94
83) Chrysene	20.98	228	92011m	3.177	ng	
88) Benzo(b)fluoranthene	22.41	252	109078m	3.497	ng	
90) Benzo(a)pyrene	22.92	252	86202	2.977	ng	# 90
92) Benzo(g,h,i)perylene	25.60	276	54905	2.095	ng	93

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

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