

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP052920\
 Data File : BP002324.D
 Acq On : 29 May 2020 20:18
 Operator : CG/JU
 Sample : L2746-05 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NM109-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/1/2020 12:24:32 PM

Quant Time: May 30 02:08:12 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP052720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 29 15:09:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	310837	20.00	ng	0.00
21) Naphthalene-d8	10.38	136	1286784	20.00	ng	0.00
39) Acenaphthene-d10	14.25	164	773320	20.00	ng	0.00
64) Phenanthrene-d10	17.01	188	1646849	20.00	ng	0.00
76) Chrysene-d12	21.11	240	1458926	20.00	ng	0.00
86) Perylene-d12	23.32	264	1527440	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	700073	42.30	ng	0.00
7) Phenol-d6	6.79	99	974463	43.28	ng	0.00
23) Nitrobenzene-d5	8.75	82	622137	29.07	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	320760	43.76	ng	0.00
45) 2-Fluorobiphenyl	12.87	172	1392138	31.77	ng	0.00
79) Terphenyl-d14	19.59	244	2019079	33.52	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.71	163	429068	8.197	ng	99
71) Phenanthrene	17.05	178	374495	4.801	ng	98
75) Fluoranthene	19.03	202	776268	8.640	ng	99
78) Pyrene	19.38	202	740213	8.148	ng	99
81) Benzo(a)anthracene	21.09	228	407720	4.882	ng	99
83) Chrysene	21.14	228	382911	4.777	ng	98
87) Indeno(1,2,3-cd)pyrene	25.53	276	258546	2.735	ng	97
88) Benzo(b)fluoranthene	22.66	252	527270	6.332	ng	# 96
89) Benzo(k)fluoranthene	22.70	252	166096m	2.078	ng	# 94
90) Benzo(a)pyrene	23.22	252	374114	4.843	ng	# 94
92) Benzo(g,h,i)perylene	26.20	276	268274	3.412	ng	# 93

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

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