

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP052920\
 Data File : BP002329.D
 Acq On : 29 May 2020 23:09
 Operator : CG/JU
 Sample : L2776-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 339

Manual Integrations
 APPROVED

mohammad
 6/1/2020 12:24:38 PM

Quant Time: May 30 10:01:04 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	319328	20.00	ng	0.00
21) Naphthalene-d8	10.38	136	1308200	20.00	ng	0.00
39) Acenaphthene-d10	14.25	164	773840	20.00	ng	0.00
64) Phenanthrene-d10	17.01	188	1592390	20.00	ng	0.00
76) Chrysene-d12	21.11	240	1353534	20.00	ng	0.00
86) Perylene-d12	23.32	264	1372387	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	2069601	121.72	ng	0.00
7) Phenol-d6	6.79	99	2855368	123.45	ng	0.00
23) Nitrobenzene-d5	8.75	82	1874890	86.17	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	904004	123.26	ng	0.00
45) 2-Fluorobiphenyl	12.87	172	3749478	85.50	ng	0.00
79) Terphenyl-d14	19.60	244	4602954	82.36	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.72	163	684231	13.063	ng	99
71) Phenanthrene	17.05	178	212205	2.813	ng	99
75) Fluoranthene	19.03	202	445876	5.132	ng	99
78) Pyrene	19.39	202	451186	5.353	ng	98
81) Benzo(a)anthracene	21.10	228	206831	2.669	ng	93
83) Chrysene	21.14	228	237520m	3.194	ng	
84) Bis(2-ethylhexyl)phthalate	21.06	149	219767	4.286	ng	98
88) Benzo(b)fluoranthene	22.66	252	301305m	4.027	ng	
90) Benzo(a)pyrene	23.22	252	197401	2.844	ng	# 94
92) Benzo(g,h,i)perylene	26.22	276	151261	2.141	ng	# 92

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

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