

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060120\
 Data File : BP002341.D
 Acq On : 01 Jun 2020 16:18
 Operator : CG/JU
 Sample : L2745-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NM81-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/2/2020 3:23:43 PM

Quant Time: Jun 03 06:27:05 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	7.61	152	293840	20.00	ng	0.00
21) Naphthalene-d8	10.38	136	1211825	20.00	ng	0.00
39) Acenaphthene-d10	14.26	164	727595	20.00	ng	0.00
64) Phenanthrene-d10	17.01	188	1540963	20.00	ng	0.00
76) Chrysene-d12	21.12	240	1332239	20.00	ng	0.00
86) Perylene-d12	23.33	264	1374163	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	2043802	130.63	ng	0.00
7) Phenol-d6	6.80	99	2861599	134.46	ng	0.00
23) Nitrobenzene-d5	8.75	82	1993097	98.88	ng	0.00
42) 2,4,6-Tribromophenol	15.76	330	987137	143.15	ng	0.00
45) 2-Fluorobiphenyl	12.87	172	4021635	97.53	ng	0.00
79) Terphenyl-d14	19.60	244	5033979	91.51	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.72	163	271236	5.507	ng	100
71) Phenanthrene	17.06	178	599014	8.207	ng	98
75) Fluoranthene	19.04	202	1266944	15.070	ng	99
78) Pyrene	19.39	202	1157880	13.957	ng	99
81) Benzo(a)anthracene	21.10	228	617732	8.100	ng	99
83) Chrysene	21.15	228	577733	7.894	ng	97
87) Indeno(1,2,3-cd)pyrene	25.56	276	405219	4.764	ng	97
88) Benzo(b)fluoranthene	22.67	252	774823m	10.342	ng	
89) Benzo(k)fluoranthene	22.70	252	308612m	4.293	ng	
90) Benzo(a)pyrene	23.23	252	584477	8.410	ng	# 97
92) Benzo(g,h,i)perylene	26.23	276	410624	5.805	ng	# 95

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

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