

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP060120\
 Data File : BP002336.D
 Acq On : 01 Jun 2020 13:29
 Operator : CG/JU
 Sample : L2773-14
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NM25-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/2/2020 3:23:32 PM

Quant Time: Jun 02 06:50:05 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	7.60	152	389300	20.00	ng	0.00
21) Naphthalene-d8	10.38	136	1579158	20.00	ng	0.00
39) Acenaphthene-d10	14.25	164	936219	20.00	ng	0.00
64) Phenanthrene-d10	17.01	188	1957070	20.00	ng	0.00
76) Chrysene-d12	21.11	240	1823836	20.00	ng	0.00
86) Perylene-d12	23.32	264	1872440	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	1336156	64.46	ng	0.00
7) Phenol-d6	6.79	99	1958090	69.44	ng	0.00
23) Nitrobenzene-d5	8.75	82	1328010	50.56	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	660923	74.49	ng	0.00
45) 2-Fluorobiphenyl	12.87	172	2783140	52.46	ng	0.00
79) Terphenyl-d14	19.60	244	3790426	50.33	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.72	163	478052	7.544	ng	99
71) Phenanthrene	17.05	178	332604	3.588	ng	99
75) Fluoranthene	19.03	202	743005	6.959	ng	99
78) Pyrene	19.39	202	710635	6.257	ng	98
81) Benzo(a)anthracene	21.10	228	399895	3.830	ng	99
83) Chrysene	21.15	228	390109m	3.893	ng	
87) Indeno(1,2,3-cd)pyrene	25.54	276	270238	2.332	ng	97
88) Benzo(b)fluoranthene	22.66	252	515510	5.050	ng	# 97
89) Benzo(k)fluoranthene	22.70	252	200026m	2.042	ng	
90) Benzo(a)pyrene	23.22	252	390438	4.123	ng	# 97
92) Benzo(a,h,i)perylene	26.22	276	286025	2.967	ng	# 94

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

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