

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP052920\
 Data File : BP002321.D
 Acq On : 29 May 2020 18:36
 Operator : CG/JU
 Sample : L2773-21 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 NM91-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/1/2020 12:24:26 PM

Quant Time: May 30 01:29:57 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	337732	20.00	ng	0.00
21) Naphthalene-d8	10.37	136	1435034	20.00	ng	0.00
39) Acenaphthene-d10	14.25	164	837237	20.00	ng	0.00
64) Phenanthrene-d10	17.00	188	1723216	20.00	ng	0.00
76) Chrysene-d12	21.11	240	1565437	20.00	ng	0.00
86) Perylene-d12	23.32	264	1582303	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	1065983	59.28	ng	0.00
7) Phenol-d6	6.79	99	1496415	61.17	ng	0.00
23) Nitrobenzene-d5	8.75	82	978263	40.99	ng	0.00
42) 2,4,6-Tribromophenol	15.75	330	477017	60.12	ng	0.00
45) 2-Fluorobiphenyl	12.87	172	2115369	44.58	ng	0.00
79) Terphenyl-d14	19.59	244	2765631	42.78	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.71	163	251701	4.441	ng	99
71) Phenanthrene	17.05	178	456873	5.597	ng	99
75) Fluoranthene	19.03	202	933608	9.931	ng	99
78) Pyrene	19.38	202	820364	8.416	ng	98
81) Benzo(a)anthracene	21.09	228	436411	4.870	ng	99
83) Chrysene	21.15	228	379339	4.411	ng	99
87) Indeno(1,2,3-cd)pyrene	25.53	276	266599	2.722	ng	96
88) Benzo(b)fluoranthene	22.66	252	519092m	6.017	ng	
89) Benzo(k)fluoranthene	22.69	252	188315m	2.275	ng	
90) Benzo(a)pyrene	23.22	252	390830	4.884	ng	# 97
92) Benzo(a,h,i)perylene	26.20	276	275543	3.383	ng	# 93

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

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