

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP061920\
 Data File : BP002467.D
 Acq On : 19 Jun 2020 19:59
 Operator : CG/JU
 Sample : L3035-13
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 2-SP-3-5

Manual Integrations
 APPROVED

mohammad
 6/22/2020 3:26:15 PM

Quant Time: Jun 20 04:17:24 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 05 15:08:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	191519	20.00	ng	0.00
21) Naphthalene-d8	10.36	136	817971	20.00	ng	0.00
39) Acenaphthene-d10	14.23	164	510127	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	1136685	20.00	ng	-0.01
76) Chrysene-d12	21.09	240	1065949	20.00	ng	-0.02
86) Perylene-d12	23.29	264	896037	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	985567	95.22	ng	0.00
7) Phenol-d6	6.78	99	1456393	105.69	ng	0.00
23) Nitrobenzene-d5	8.73	82	1081208	78.94	ng	-0.01
42) 2,4,6-Tribromophenol	15.73	330	526601	107.82	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	2221163	73.21	ng	0.00
79) Terphenyl-d14	19.58	244	3340187	78.38	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.69	163	334679	9.575	ng	100
71) Phenanthrene	17.03	178	376987	7.090	ng	100
72) Anthracene	17.12	178	113244	2.144	ng	98
75) Fluoranthene	19.02	202	779525	12.282	ng	98
78) Pyrene	19.37	202	704049	10.783	ng	97
81) Benzo(a)anthracene	21.08	228	368559	6.057	ng	98
83) Chrysene	21.13	228	314546m	5.457	ng	
87) Indeno(1,2,3-cd)pyrene	25.49	276	166792	2.976	ng	97
88) Benzo(b)fluoranthene	22.63	252	371251	7.683	ng	# 96
89) Benzo(k)fluoranthene	22.67	252	118048m	2.571	ng	
90) Benzo(a)pyrene	23.19	252	268829	5.937	ng	97
92) Benzo(a,h,i)perylene	26.16	276	159141	3.418	ng	97

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

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