

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP061920\
 Data File : BP002460.D
 Acq On : 19 Jun 2020 16:01
 Operator : CG/JU
 Sample : L3055-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 AMITY-ST-ENTRANCE

Manual Integrations
 APPROVED

mohammad
 6/22/2020 3:26:10 PM

Quant Time: Jun 19 16:42:00 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	183581	20.00	ng	0.00
21) Naphthalene-d8	10.36	136	682631	20.00	ng	0.00
39) Acenaphthene-d10	14.23	164	367430	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	739442	20.00	ng	0.00
76) Chrysene-d12	21.10	240	683546	20.00	ng	-0.01
86) Perylene-d12	23.30	264	717100	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	1175565	118.49	ng	0.00
7) Phenol-d6	6.78	99	1549979	117.34	ng	0.00
23) Nitrobenzene-d5	8.73	82	1041068	91.08	ng	-0.01
42) 2,4,6-Tribromophenol	15.73	330	436943	124.21	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	1924407	88.06	ng	0.00
79) Terphenyl-d14	19.58	244	2429917	94.46	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.69	163	237339	9.427	ng	99
75) Fluoranthene	19.02	202	397910	9.637	ng	98
78) Pyrene	19.37	202	441349	10.541	ng	98
81) Benzo(a)anthracene	21.08	228	512506	13.136	ng	97
83) Chrysene	21.13	228	499573m	13.515	ng	
87) Indeno(1,2,3-cd)pyrene	25.51	276	267046	5.954	ng	97
88) Benzo(b)fluoranthene	22.64	252	460859m	11.917	ng	
89) Benzo(k)fluoranthene	22.68	252	132629m	3.610	ng	
90) Benzo(a)pyrene	23.20	252	611202	16.866	ng	96
91) Dibenzo(a,h)anthracene	25.52	278	209852	5.833	ng	# 75
92) Benzo(g,h,i)perylene	26.19	276	400792	10.756	ng	97

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

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