

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070120\
 Data File : BP002613.D
 Acq On : 01 Jul 2020 20:12
 Operator : CG/JU
 Sample : L3162-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GL-SP

Manual Integrations
 APPROVED

mohammad
 7/7/2020 8:23:15 AM

Quant Time: Jul 02 08:01:14 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 29 16:51:57 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	341255	20.00	ng	0.00
21) Naphthalene-d8	10.34	136	1326385	20.00	ng	0.00
39) Acenaphthene-d10	14.22	164	775118	20.00	ng	0.00
64) Phenanthrene-d10	16.97	188	1464451	20.00	ng	0.00
76) Chrysene-d12	21.09	240	1014229	20.00	ng	0.00
86) Perylene-d12	23.29	264	949375	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	1465688	72.72	ng	0.00
7) Phenol-d6	6.76	99	1997072	70.99	ng	0.00
23) Nitrobenzene-d5	8.72	82	1317334	51.17	ng	0.00
42) 2,4,6-Tribromophenol	15.72	330	615001	65.23	ng	0.00
45) 2-Fluorobiphenyl	12.83	172	2653707	50.39	ng	0.00
79) Terphenyl-d14	19.57	244	2875228	61.19	ng	0.00
Target Compounds						
49) Acenaphthylene	13.93	152	249919	3.327	ng	98
50) Dimethylphthalate	13.68	163	311343	5.114	ng	100
70) Pentachlorophenol	16.65	266	1149	2.772	ng	# 84
71) Phenanthrene	17.02	178	299875	3.806	ng	99
75) Fluoranthene	19.00	202	602483	6.317	ng	97
78) Pyrene	19.36	202	602925	8.734	ng	100
81) Benzo(a)anthracene	21.07	228	337444	5.089	ng	96
83) Chrysene	21.12	228	339596	5.341	ng	97
87) Indeno(1,2,3-cd)pyrene	25.49	276	261192	3.821	ng	99
88) Benzo(b)fluoranthene	22.63	252	470056m	7.759	ng	
89) Benzo(k)fluoranthene	22.67	252	142266m	2.517	ng	
90) Benzo(a)pyrene	23.19	252	315204	5.670	ng	# 94
92) Benzo(a,h,i)perylene	26.16	276	277650	4.970	ng	96

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

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