

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP080520\
 Data File : BP002907.D
 Acq On : 05 Aug 2020 11:48
 Operator : CG/JU
 Sample : L3541-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TL-7

Manual Integrations
 APPROVED

mohammad
 8/6/2020 1:21:15 PM

Quant Time: Aug 05 12:20:43 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 30 15:30:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	409675	20.00	ng	0.00
21) Naphthalene-d8	10.49	136	1594221	20.00	ng	0.00
39) Acenaphthene-d10	14.35	164	990194	20.00	ng	0.00
64) Phenanthrene-d10	17.10	188	2142309	20.00	ng	-0.01
76) Chrysene-d12	21.19	240	2467661	20.00	ng	0.00
86) Perylene-d12	23.46	264	2503612	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	2087454	77.98	ng	0.00
7) Phenol-d6	6.89	99	2649415	71.29	ng	0.00
23) Nitrobenzene-d5	8.86	82	1557298	52.91	ng	0.00
42) 2,4,6-Tribromophenol	15.84	330	1139948	92.63	ng	0.00
45) 2-Fluorobiphenyl	12.98	172	3828545	52.83	ng	0.00
79) Terphenyl-d14	19.67	244	5548489	45.01	ng	-0.01
Target Compounds						
50) Dimethylphthalate	13.81	163	451444	5.829	ng	99
71) Phenanthrene	17.15	178	1418463	11.291	ng	98
72) Anthracene	17.23	178	408893	3.330	ng	99
75) Fluoranthene	19.12	202	2829774	19.551	ng	99
78) Pyrene	19.47	202	2414983	13.999	ng	98
80) Butylbenzylphthalate	20.36	149	9596	3.186	ng	95
81) Benzo(a)anthracene	21.18	228	1698330	10.104	ng	98
83) Chrysene	21.23	228	1396061m	8.805	ng	
87) Indeno(1,2,3-cd)pyrene	25.76	276	785032	3.915	ng	96
88) Benzo(b)fluoranthene	22.78	252	1789264	10.127	ng	100
89) Benzo(k)fluoranthene	22.82	252	730119m	4.372	ng	
90) Benzo(a)pyrene	23.36	252	1314883	8.237	ng	98
92) Benzo(a,h,i)perylene	26.46	276	728355	4.414	ng	99

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

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