

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP061920\
 Data File : BP002465.D
 Acq On : 19 Jun 2020 18:51
 Operator : CG/JU
 Sample : L3052-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 IWS

Manual Integrations
 APPROVED

mohammad
 6/22/2020 3:26:12 PM

Quant Time: Jun 20 02:42:29 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	208513	20.00	ng	0.00
21) Naphthalene-d8	10.36	136	869779	20.00	ng	0.00
39) Acenaphthene-d10	14.23	164	536375	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	1163008	20.00	ng	0.00
76) Chrysene-d12	21.09	240	906309	20.00	ng	-0.02
86) Perylene-d12	23.29	264	810997	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	983996	87.32	ng	0.00
7) Phenol-d6	6.78	99	1374145	91.59	ng	0.00
23) Nitrobenzene-d5	8.73	82	944378	64.84	ng	-0.01
42) 2,4,6-Tribromophenol	15.73	330	511811	99.66	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	1965509	61.61	ng	0.00
79) Terphenyl-d14	19.58	244	2814110	77.29	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.69	163	257703	7.012	ng	99
71) Phenanthrene	17.03	178	129114	2.373	ng	99
75) Fluoranthene	19.02	202	278992	4.296	ng	98
78) Pyrene	19.37	202	287728	5.183	ng	98
81) Benzo(a)anthracene	21.08	228	118778	2.296	ng	95
83) Chrysene	21.13	228	138805	2.832	ng	96
88) Benzo(b)fluoranthene	22.63	252	146695m	3.354	ng	
90) Benzo(a)pyrene	23.19	252	88340	2.155	ng	# 96

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

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