

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP072220\
 Data File : BP002827.D
 Acq On : 22 Jul 2020 16:23
 Operator : CG/JU
 Sample : L3365-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20017-20405

Quant Time: Jul 23 02:23:44 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 10:54:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	144448	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	581264	20.00	ng	-0.01
39) Acenaphthene-d10	14.20	164	356102	20.00	ng	-0.01
64) Phenanthrene-d10	16.96	188	749089	20.00	ng	-0.01
76) Chrysene-d12	21.07	240	672680	20.00	ng	0.00
86) Perylene-d12	23.26	264	592151	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.18	112	849670	99.24	ng	0.00
7) Phenol-d6	6.75	99	1189858	97.38	ng	0.00
23) Nitrobenzene-d5	8.69	82	823697	75.79	ng	-0.01
42) 2,4,6-Tribromophenol	15.70	330	392594	105.62	ng	0.00
45) 2-Fluorobiphenyl	12.81	172	1646915	71.55	ng	-0.01
79) Terphenyl-d14	19.55	244	1856756	63.28	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.66	163	222056	8.177	ng	99
71) Phenanthrene	17.00	178	88815	2.190	ng	# 96
75) Fluoranthene	18.99	202	310464	6.606	ng	99
78) Pyrene	19.34	202	260401	5.607	ng	99
81) Benzo(a)anthracene	21.06	228	123052	2.826	ng	98
83) Chrysene	21.10	228	169455	4.088	ng	98
87) Indeno(1,2,3-cd)pyrene	25.47	276	90630	2.118	ng	96
88) Benzo(b)fluoranthene	22.61	252	227731	6.194	ng	# 97
90) Benzo(a)pyrene	23.17	252	121243	3.480	ng	# 97
92) Benzo(g,h,i)perylene	26.14	276	89550	2.559	ng	97

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

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