

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051520\
 Data File : BF120335.D
 Acq On : 15 May 2020 21:35
 Operator : MA/SJ
 Sample : L2505-01 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-02-051220

Quant Time: May 16 05:23:39 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 15 12:31:31 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.61	152	238791	20.00	ng	0.00
21) Naphthalene-d8	7.90	136	1020457	20.00	ng	0.00
39) Acenaphthene-d10	9.65	164	460892	20.00	ng	0.00
64) Phenanthrene-d10	11.13	188	697223	20.00	ng	0.00
76) Chrysene-d12	13.77	240	436035	20.00	ng	0.00
87) Perylene-d12	15.16	264	379573	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	374682	30.53	ng	0.00
7) Phenol-d6	6.27	99	480606	30.97	ng	0.00
23) Nitrobenzene-d5	7.17	82	303537	20.37	ng	-0.02
42) 2,4,6-Tribromophenol	10.44	330	105784	28.15	ng	0.00
45) 2-Fluorobiphenyl	8.97	172	564928	25.18	ng	0.00
79) Terphenyl-d14	12.72	244	495139	25.09	ng	0.00

Target Compounds				Qvalue		
50) Dimethylphthalate	9.37	163	79885	2.742	ng	99
58) Fluorene	10.20	166	80706	3.318	ng	95
71) Phenanthrene	11.16	178	571373	19.503	ng	98
72) Anthracene	11.20	178	246128	7.847	ng	97
75) Fluoranthene	12.35	202	1577317	50.230	ng	97
78) Pyrene	12.57	202	1285501	41.354	ng	98
81) Benzo(a)anthracene	13.76	228	601529	26.156	ng	98
83) Chrysene	13.80	228	537626	21.770	ng	96
86) Indeno(1,2,3-cd)pyrene	16.49	276	328007m	14.634	ng	
88) Benzo(b)fluoranthene	14.77	252	737229m	35.564	ng	
89) Benzo(k)fluoranthene	14.80	252	215770m	12.014	ng	
90) Benzo(a)pyrene	15.10	252	432061	23.241	ng	97
91) Dibenzo(a,h)anthracene	16.49	278	85259	5.177	ng	# 78
92) Benzo(g,h,i)perylene	16.88	276	256574	15.449	ng	96

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

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