

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031720\
 Data File : BG044840.D
 Acq On : 17 Mar 2020 19:50
 Operator : CG/JU
 Sample : L1912-09
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 Z3-S2

Quant Time: Mar 18 01:57:32 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG031020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 16 17:37:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	108105	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	447824	20.00	ng	0.00
39) Acenaphthene-d10	14.68	164	289348	20.00	ng	0.00
64) Phenanthrene-d10	17.42	188	589091	20.00	ng	0.00
76) Chrysene-d12	21.69	240	501792	20.00	ng	0.00
87) Perylene-d12	24.90	264	564335	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	673603	97.00	ng	0.00
7) Phenol-d6	7.20	99	974107	96.54	ng	0.00
23) Nitrobenzene-d5	9.20	82	631657	61.89	ng	0.00
42) 2,4,6-Tribromophenol	16.16	330	368747	97.33	ng	0.00
45) 2-Fluorobiphenyl	13.30	172	1287499	63.72	ng	0.00
79) Terphenyl-d14	20.04	244	1661349	61.93	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.13	163	149423	6.484	ng	99
71) Phenanthrene	17.46	178	113190	3.596	ng	99
75) Fluoranthene	19.47	202	232839	6.212	ng	95
78) Pyrene	19.83	202	217582	5.910	ng	98
81) Benzo(a)anthracene	21.67	228	121689	3.368	ng	97
83) Chrysene	21.74	228	111979	3.244	ng	96
88) Benzo(b)fluoranthene	23.88	252	148432	3.984	ng	97
90) Benzo(a)pyrene	24.75	252	116138	3.331	ng	99
92) Benzo(g,h,i)perylene	29.67	276	73177	2.106	ng	# 92

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

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