

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG041620\
 Data File : BG045016.D
 Acq On : 16 Apr 2020 21:30
 Operator : CG/JU
 Sample : L2230-29
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-20

Quant Time: Apr 17 05:29:50 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 15 17:57:00 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	109231	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	471135	20.00	ng	0.00
39) Acenaphthene-d10	14.65	164	336608	20.00	ng	0.00
64) Phenanthrene-d10	17.40	188	728843	20.00	ng	0.00
76) Chrysene-d12	21.66	240	654531	20.00	ng	0.00
87) Perylene-d12	24.85	264	732083	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	379558	57.37	ng	0.00
7) Phenol-d6	7.18	99	580587	59.06	ng	0.00
23) Nitrobenzene-d5	9.18	82	426626	41.32	ng	0.00
42) 2,4,6-Tribromophenol	16.14	330	302668	58.20	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	1032681	44.99	ng	0.00
79) Terphenyl-d14	20.01	244	1428310	47.56	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.10	163	167469	6.110	ng	Qvalue # 98
71) Phenanthrene	17.44	178	107761	2.877	ng	95
78) Pyrene	19.81	202	269999	5.984	ng	94
81) Benzo(a)anthracene	21.64	228	185055	4.362	ng	95
83) Chrysene	21.71	228	177646	4.358	ng	98
86) Indeno(1,2,3-cd)pyrene	28.46	276	137079	2.533	ng	# 93
88) Benzo(b)fluoranthene	23.84	252	245703	5.358	ng	98
90) Benzo(a)pyrene	24.69	252	186304	4.303	ng	99
92) Benzo(g,h,i)perylene	29.56	276	131828	3.014	ng	99

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

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