

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101519\
 Data File : BG043229.D
 Acq On : 15 Oct 2019 21:44
 Operator : JU
 Sample : K5148-05 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OR-03-101119

Manual Integrations
 APPROVED

mohammad
 10/17/2019 7:57:27 AM

Quant Time: Oct 16 03:41:45 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 09 01:11:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	65796	20.00	ng	0.00
21) Naphthalene-d8	10.86	136	250646	20.00	ng	0.00
39) Acenaphthene-d10	14.67	164	175303	20.00	ng	0.00
64) Phenanthrene-d10	17.42	188	371690	20.00	ng	0.00
76) Chrysene-d12	21.70	240	404370	20.00	ng	0.00
87) Perylene-d12	24.93	264	484635	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	163063	44.23	ng	0.00
7) Phenol-d6	7.22	99	238567	44.93	ng	0.00
23) Nitrobenzene-d5	9.22	82	142552	27.64	ng	0.00
42) 2,4,6-Tribromophenol	16.16	330	134611	44.66	ng	0.00
45) 2-Fluorobiphenyl	13.30	172	388746	32.15	ng	0.00
79) Terphenyl-d14	20.03	244	661251	34.85	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.12	163	60990	4.643	ng	Qvalue # 97
71) Phenanthrene	17.47	178	86364	4.759	ng	98
75) Fluoranthene	19.48	202	212816	8.867	ng	96
78) Pyrene	19.83	202	212173	8.791	ng	99
81) Benzo(a)anthracene	21.67	228	114040	4.526	ng	96
83) Chrysene	21.74	228	120375	4.923	ng	96
86) Indeno(1,2,3-cd)pyrene	28.59	276	83515	2.744	ng	# 92
88) Benzo(b)fluoranthene	23.91	252	163354	5.957	ng	# 93
89) Benzo(k)fluoranthene	23.95	252	54591m	2.012	ng	
90) Benzo(a)pyrene	24.77	252	124390	4.808	ng	# 95
92) Benzo(g,h,i)perylene	29.75	276	89032	3.464	ng	98

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

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