

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110619\
 Data File : BG043534.D
 Acq On : 7 Nov 2019 00:01
 Operator : JU
 Sample : K5675-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TR-05-110519

Manual Integrations
 APPROVED

Sohil
 11/8/2019 3:29:58 PM

Quant Time: Nov 07 02:34:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 04 15:22:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	41241	20.00	ng	0.00
21) Naphthalene-d8	10.81	136	149031	20.00	ng	0.00
39) Acenaphthene-d10	14.64	164	100427	20.00	ng	0.00
64) Phenanthrene-d10	17.38	188	223004	20.00	ng	0.00
76) Chrysene-d12	21.65	240	228514	20.00	ng	0.00
87) Perylene-d12	24.84	264	269379	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	161348	71.69	ng	0.00
7) Phenol-d6	7.19	99	229067	70.74	ng	0.00
23) Nitrobenzene-d5	9.17	82	137285	47.49	ng	0.00
42) 2,4,6-Tribromophenol	16.13	330	109193	57.14	ng	0.00
45) 2-Fluorobiphenyl	13.26	172	351036	48.30	ng	0.00
79) Terphenyl-d14	20.00	244	605876	49.74	ng	0.00
Target Compounds						
10) Phenol	7.22	94	6952	2.349	ng	79
50) Dimethylphthalate	14.09	163	64363	8.274	ng	# 95
71) Phenanthrene	17.43	178	36484	3.195	ng	98
75) Fluoranthene	19.44	202	177223	11.904	ng	97
78) Pyrene	19.80	202	167788	11.862	ng	98
81) Benzo(a)anthracene	21.63	228	94823	6.625	ng	97
83) Chrysene	21.70	228	87636	6.162	ng	97
86) Indeno(1,2,3-cd)pyrene	28.48	276	75577	4.233	ng	# 96
88) Benzo(b)fluoranthene	23.84	252	140077	9.181	ng	# 94
89) Benzo(k)fluoranthene	23.89	252	46220m	3.009	ng	
90) Benzo(a)pyrene	24.70	252	104616	7.181	ng	99
92) Benzo(a,h,i)perylene	29.60	276	71867m	4.889	ng	

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

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