

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110519\
 Data File : BG043514.D
 Acq On : 5 Nov 2019 19:07
 Operator : JU
 Sample : K5602-05RX
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20190942-AMENDED-COMPOSITE-

Manual Integrations
 APPROVED

mohammad
 11/6/2019 8:41:21 AM

Quant Time: Nov 06 02:27:02 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	31876	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	117376	20.00	ng	0.00
39) Acenaphthene-d10	14.64	164	89272	20.00	ng	0.00
64) Phenanthrene-d10	17.38	188	227451	20.00	ng	0.00
76) Chrysene-d12	21.66	240	247194	20.00	ng	0.00
87) Perylene-d12	24.85	264	294702	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	5772	3.32	ng	0.00
7) Phenol-d6	7.20	99	29936	11.96	ng	0.00
23) Nitrobenzene-d5	9.17	82	141312	62.07	ng	0.00
42) 2,4,6-Tribromophenol	16.15	330	6162	3.63	ng	0.01
45) 2-Fluorobiphenyl	13.26	172	416778	64.51	ng	0.00
79) Terphenyl-d14	19.99	244	861875	65.41	ng	0.00
Target Compounds						
10) Phenol	7.23	94	6608	2.889	ng	99
50) Dimethylphthalate	14.09	163	71187	10.295	ng	97
75) Fluoranthene	19.44	202	51055	3.362	ng	99
78) Pyrene	19.80	202	48578	3.175	ng	98
81) Benzo(a)anthracene	21.63	228	31581	2.040	ng	95
83) Chrysene	21.70	228	34814	2.263	ng	95
88) Benzo(b)fluoranthene	23.84	252	39144	2.345	ng	99

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

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