

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090319\
 Data File : BG042697.D
 Acq On : 3 Sep 2019 20:24
 Operator : HP/JU
 Sample : K4602-08
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PAR-C1-C4-(0-3.11)

Manual Integrations
 APPROVED

mohammad
 9/6/2019 8:17:40 AM

Quant Time: Sep 04 04:58:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 22 14:29:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	34561	20.00	ng	-0.01
21) Naphthalene-d8	10.82	136	135473	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	105295	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	232034	20.00	ng	-0.01
76) Chrysene-d12	21.68	240	257404	20.00	ng	-0.01
87) Perylene-d12	24.92	264	309640	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	141369	82.84	ng	0.00
7) Phenol-d6	7.16	99	191618	83.13	ng	-0.01
23) Nitrobenzene-d5	9.16	82	112277	57.27	ng	-0.01
42) 2,4,6-Tribromophenol	16.15	330	132404	88.47	ng	-0.01
45) 2-Fluorobiphenyl	13.28	172	315627	50.70	ng	0.00
79) Terphenyl-d14	20.02	244	549684	46.17	ng	0.00
Target Compounds						
10) Phenol	7.19	94	9747	4.159	ng	88
50) Dimethylphthalate	14.10	163	43334	5.704	ng	98
75) Fluoranthene	19.46	202	37631	2.675	ng	93
78) Pyrene	19.83	202	34984m	2.217	ng	
83) Chrysene	21.73	228	35475	2.349	ng	94
88) Benzo(b)fluoranthene	23.90	252	59839	3.515	ng	99
90) Benzo(a)pyrene	24.77	252	35283	2.184	ng	98

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

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