

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091620\
 Data File : BG046647.D
 Acq On : 16 Sep 2020 19:21
 Operator : CG/JU
 Sample : L4021-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B14-5-7

Quant Time: Sep 17 00:52:16 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 16 11:07:52 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	68398	20.00	ng	0.00
21) Naphthalene-d8	10.72	136	283398	20.00	ng	0.00
39) Acenaphthene-d10	14.55	164	188397	20.00	ng	0.00
64) Phenanthrene-d10	17.29	188	427513	20.00	ng	0.00
76) Chrysene-d12	21.54	240	429674	20.00	ng	0.00
86) Perylene-d12	24.64	264	429454	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	332888	86.20	ng	0.00
7) Phenol-d6	7.09	99	473112	86.42	ng	0.00
23) Nitrobenzene-d5	9.07	82	310127	62.55	ng	0.00
42) 2,4,6-Tribromophenol	16.04	330	189284	74.15	ng	0.00
45) 2-Fluorobiphenyl	13.17	172	831065	67.34	ng	0.00
79) Terphenyl-d14	19.91	244	1213171	60.09	ng	0.00
Target Compounds						Qvalue
10) Phenol	7.12	94	24418	4.843	ng	85
50) Dimethylphthalate	14.00	163	98993	6.792	ng	99
71) Phenanthrene	17.34	178	57935	2.674	ng	97
75) Fluoranthene	19.35	202	124369	4.460	ng	97
78) Pyrene	19.71	202	114342	4.179	ng	97
81) Benzo(a)anthracene	21.52	228	62898	2.349	ng	95
83) Chrysene	21.58	228	65207	2.531	ng	96
88) Benzo(b)fluoranthene	23.66	252	85178	3.176	ng	98
90) Benzo(a)pyrene	24.49	252	57120	2.283	ng	98

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

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