

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091620\
 Data File : BG046646.D
 Acq On : 16 Sep 2020 18:41
 Operator : CG/JU
 Sample : L4021-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B15-1-3

Quant Time: Sep 17 00:48:29 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	72049	20.00	ng	0.00
21) Naphthalene-d8	10.72	136	289791	20.00	ng	0.00
39) Acenaphthene-d10	14.55	164	184288	20.00	ng	0.00
64) Phenanthrene-d10	17.30	188	421650	20.00	ng	0.00
76) Chrysene-d12	21.54	240	434650	20.00	ng	0.00
86) Perylene-d12	24.63	264	429926	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	371921	91.43	ng	0.00
7) Phenol-d6	7.10	99	493906	85.65	ng	0.00
23) Nitrobenzene-d5	9.08	82	328273	64.74	ng	0.00
42) 2,4,6-Tribromophenol	16.04	330	202490	81.09	ng	0.00
45) 2-Fluorobiphenyl	13.17	172	794462	65.81	ng	0.00
79) Terphenyl-d14	19.91	244	1153432	56.48	ng	0.00
Target Compounds						
10) Phenol	7.12	94	25711	4.841	ng	91
50) Dimethylphthalate	14.00	163	92050	6.456	ng	98
71) Phenanthrene	17.34	178	56301	2.634	ng	95
75) Fluoranthene	19.35	202	110227	4.008	ng	96
78) Pyrene	19.71	202	93100	3.364	ng	95
88) Benzo(b)fluoranthene	23.67	252	59350	2.211	ng	95

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

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