

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG041520\
 Data File : BG045001.D
 Acq On : 16 Apr 2020 3:34
 Operator : CG/JU
 Sample : L2278-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CHASSIS-WASH

Quant Time: Apr 16 05:42:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 15 17:57:00 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	65097	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	277674	20.00	ng	0.00
39) Acenaphthene-d10	14.65	164	221251	20.00	ng	0.00
64) Phenanthrene-d10	17.40	188	534654	20.00	ng	0.00
76) Chrysene-d12	21.66	240	504610	20.00	ng	0.00
87) Perylene-d12	24.85	264	550464	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	465802	118.13	ng	0.00
7) Phenol-d6	7.18	99	685766	117.06	ng	0.00
23) Nitrobenzene-d5	9.18	82	536838	88.22	ng	0.00
42) 2,4,6-Tribromophenol	16.15	330	521072	152.45	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	1371796	90.91	ng	0.00
79) Terphenyl-d14	20.01	244	2390092	103.24	ng	0.00
Target Compounds						
10) Phenol	7.21	94	20189	3.444	ng	Qvalue # 71
15) Benzyl Alcohol	8.26	79	186029	37.875	ng	92
17) 2-Methylphenol	8.46	107	9316	2.060	ng	90
20) 3+4-Methylphenols	8.79	107	67632	10.683	ng	93
32) Benzoic acid	10.13	122	110321	34.152	ng	94

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

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