

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111119\
 Data File : BG043601.D
 Acq On : 12 Nov 2019 2:26
 Operator : JU
 Sample : K5782-06
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-2A-COMP

Quant Time: Nov 12 08:15:35 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	8.00	152	30528	20.00	ng	0.00
21) Naphthalene-d8	10.81	136	121216	20.00	ng	0.00
39) Acenaphthene-d10	14.63	164	96730	20.00	ng	0.00
64) Phenanthrene-d10	17.38	188	238258	20.00	ng	-0.01
76) Chrysene-d12	21.64	240	196696	20.00	ng	-0.01
87) Perylene-d12	24.83	264	234492	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	163616	98.21	ng	0.00
7) Phenol-d6	7.20	99	157750	65.81	ng	0.00
23) Nitrobenzene-d5	9.16	82	232975	99.09	ng	-0.01
42) 2,4,6-Tribromophenol	16.13	330	276242	150.07	ng	0.00
45) 2-Fluorobiphenyl	13.25	172	664848	94.97	ng	-0.01
79) Terphenyl-d14	19.99	244	1296559	123.66	ng	0.00
Target Compounds						
10) Phenol	7.22	94	6901	3.151	ng	Qvalue # 61
31) Naphthalene	10.86	128	88321	14.354	ng	96
37) 2-Methylnaphthalene	12.46	142	41959	8.888	ng	96
38) 1-Methylnaphthalene	12.68	142	34090	7.589	ng	99
50) Dimethylphthalate	14.08	163	59243	7.907	ng	97
52) Acenaphthene	14.70	154	18022	3.391	ng	91

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

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