

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG103119\
 Data File : BG043436.D
 Acq On : 31 Oct 2019 16:27
 Operator : JU
 Sample : K5497-05
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 169-32-(0-1)

Quant Time: Nov 01 04:58:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 21 16:37:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	43548	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	157403	20.00	ng	0.00
39) Acenaphthene-d10	14.65	164	118040	20.00	ng	0.00
64) Phenanthrene-d10	17.39	188	287486	20.00	ng	0.00
76) Chrysene-d12	21.66	240	329305	20.00	ng	0.00
87) Perylene-d12	24.86	264	390517	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	271026	114.04	ng	0.00
7) Phenol-d6	7.21	99	391387	114.47	ng	0.00
23) Nitrobenzene-d5	9.19	82	233704	76.55	ng	0.00
42) 2,4,6-Tribromophenol	16.14	330	248225	110.50	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	548436	64.20	ng	0.00
79) Terphenyl-d14	20.00	244	1047206	59.66	ng	0.00
Target Compounds						
10) Phenol	7.24	94	8711	2.788	ng	81
50) Dimethylphthalate	14.10	163	60168	6.581	ng	99
75) Fluoranthene	19.45	202	44394	2.313	ng	96
78) Pyrene	19.81	202	41711	2.046	ng	96

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

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