

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102819\
 Data File : BG043363.D
 Acq On : 28 Oct 2019 23:26
 Operator : JU
 Sample : K5539-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 162-36-(0-1)

Quant Time: Oct 29 13:46:55 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	44350	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	160760	20.00	ng	0.00
39) Acenaphthene-d10	14.65	164	120612	20.00	ng	0.00
64) Phenanthrene-d10	17.39	188	286395	20.00	ng	0.00
76) Chrysene-d12	21.66	240	328681	20.00	ng	0.00
87) Perylene-d12	24.86	264	381788	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	333845	137.94	ng	0.00
7) Phenol-d6	7.21	99	470947	135.24	ng	0.00
23) Nitrobenzene-d5	9.19	82	280711	90.02	ng	0.00
42) 2,4,6-Tribromophenol	16.14	330	289653	126.20	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	694934	79.62	ng	0.00
79) Terphenyl-d14	20.00	244	1336153	76.27	ng	0.00
Target Compounds						
10) Phenol	7.24	94	12550	3.944	ng	Qvalue # 75
50) Dimethylphthalate	14.10	163	60475	6.473	ng	98
71) Phenanthrene	17.44	178	53446	3.644	ng	97
75) Fluoranthene	19.45	202	95122	4.975	ng	97
78) Pyrene	19.81	202	83497	4.104	ng	96
81) Benzo(a)anthracene	21.64	228	46718	2.269	ng	97
88) Benzo(b)fluoranthene	23.84	252	44973	2.080	ng	# 96

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

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