

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082519\
 Data File : BG042480.D
 Acq On : 25 Aug 2019 16:25
 Operator : HP/JU
 Sample : K4493-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 T9-12-13-I-(0-1.9)

Quant Time: Aug 26 08:03:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 22 14:29:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	47883	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	188817	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	140184	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	341959	20.00	ng	0.00
76) Chrysene-d12	21.69	240	349086	20.00	ng	0.00
87) Perylene-d12	24.93	264	409234	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	240547	101.74	ng	0.00
7) Phenol-d6	7.16	99	332173	104.01	ng	0.00
23) Nitrobenzene-d5	9.17	82	187398	68.58	ng	0.00
42) 2,4,6-Tribromophenol	16.15	330	206256	103.52	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	555535	67.02	ng	0.00
79) Terphenyl-d14	20.02	244	991715	61.42	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	14.11	163	95525	9.444	ng	98
71) Phenanthrene	17.46	178	51364	3.024	ng	97
75) Fluoranthene	19.47	202	113968	5.498	ng	98
78) Pyrene	19.83	202	93794	4.382	ng	96
81) Benzo(a)anthracene	21.67	228	44763	2.108	ng	98
83) Chrysene	21.73	228	50842	2.483	ng	98
88) Benzo(b)fluoranthene	23.91	252	74214	3.299	ng	# 98
90) Benzo(a)pyrene	24.78	252	52070	2.439	ng	96

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

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