

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG063020\
 Data File : BG045897.D
 Acq On : 30 Jun 2020 23:28
 Operator : CG/JU
 Sample : L2819-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BU-03-062620

Manual Integrations
 APPROVED

Sohil
 7/7/2020 8:17:14 AM

Quant Time: Jul 01 03:34:42 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 30 16:34:40 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	67880	20.00	ng	0.00
21) Naphthalene-d8	10.70	136	263263	20.00	ng	0.00
39) Acenaphthene-d10	14.55	164	175281	20.00	ng	0.00
64) Phenanthrene-d10	17.30	188	351678	20.00	ng	0.00
76) Chrysene-d12	21.56	240	312955	20.00	ng	0.00
86) Perylene-d12	24.67	264	326162	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	388075	96.57	ng	0.00
7) Phenol-d6	7.12	99	559332	91.51	ng	0.00
23) Nitrobenzene-d5	9.06	82	378382	65.66	ng	0.00
42) 2,4,6-Tribromophenol	16.05	330	225865	85.35	ng	0.00
45) 2-Fluorobiphenyl	13.17	172	845573	70.89	ng	0.00
79) Terphenyl-d14	19.92	244	1124459	72.82	ng	0.00
Target Compounds						
32) Benzoic acid	10.09	122	6936m	11.202	ng	Qvalue
50) Dimethylphthalate	14.00	163	82617	6.017	ng	97
75) Fluoranthene	19.35	202	82873	3.716	ng	98
78) Pyrene	19.72	202	73292	3.497	ng	96
81) Benzo(a)anthracene	21.54	228	42469	2.093	ng	96
83) Chrysene	21.60	228	46447m	2.367	ng	
88) Benzo(b)fluoranthene	23.70	252	66109	3.233	ng	# 90
90) Benzo(a)pyrene	24.52	252	42683	2.234	ng	# 92

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

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