

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021920\
 Data File : BG044513.D
 Acq On : 20 Feb 2020 7:51
 Operator : CG/JU
 Sample : L1573-02
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Feb 20 10:26:58 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 10 16:11:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	64218	20.00	ng	-0.02
21) Naphthalene-d8	10.69	136	271601	20.00	ng	-0.01
39) Acenaphthene-d10	14.53	164	186075	20.00	ng	-0.01
64) Phenanthrene-d10	17.28	188	408324	20.00	ng	-0.01
76) Chrysene-d12	21.52	240	388472	20.00	ng	-0.02
87) Perylene-d12	24.60	264	424431	20.00	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.47	112	402697	110.67	ng	-0.01
7) Phenol-d6	7.07	99	530771	108.62	ng	-0.01
23) Nitrobenzene-d5	9.05	82	330527	68.70	ng	-0.02
42) 2,4,6-Tribromophenol	16.02	330	226412	103.65	ng	-0.01
45) 2-Fluorobiphenyl	13.15	172	790028	71.10	ng	-0.01
79) Terphenyl-d14	19.89	244	1241162	65.72	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	7.09	94	10993	2.273	ng	# 73
31) Naphthalene	10.74	128	344593	26.151	ng	98
37) 2-Methylnaphthalene	12.35	142	85574	8.747	ng	96
38) 1-Methylnaphthalene	12.57	142	41219	4.469	ng	97
46) 1,1'-Biphenyl	13.36	154	36905	2.750	ng	94
50) Dimethylphthalate	13.98	163	31973	2.390	ng	98
52) Acenaphthene	14.59	154	213228	21.000	ng	98
55) Dibenzofuran	14.93	168	229511	14.929	ng	98
58) Fluorene	15.58	166	206107	17.022	ng	99
71) Phenanthrene	17.32	178	1486836	75.197	ng	97
72) Anthracene	17.41	178	370932	18.975	ng	96
73) Carbazole	17.68	167	187525	10.406	ng	96
75) Fluoranthene	19.33	202	1421344	62.141	ng	98
78) Pyrene	19.69	202	1156893	46.373	ng	99
81) Benzo(a)anthracene	21.50	228	629858	26.308	ng	97
83) Chrysene	21.57	228	517112	22.345	ng	97
86) Indeno(1,2,3-cd)pyrene	28.08	276	290849	9.633	ng	# 93
88) Benzo(b)fluoranthene	23.63	252	683779	27.958	ng	98
89) Benzo(k)fluoranthene	23.69	252	255919m	10.736	ng	
90) Benzo(a)pyrene	24.45	252	498138	21.542	ng	99
91) Dibenzo(a,h)anthracene	28.12	278	71216m	3.112	ng	
92) Benzo(g,h,i)perylene	29.17	276	275646	12.032	ng	97

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

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