

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121522\
 Data File : BG055988.D
 Acq On : 15 Dec 2022 22:37
 Operator : CG/JU
 Sample : N6037-02 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RB-17-(1-3)

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/16/2022
 Supervised By :mohammad ahmed 12/17/2022

Quant Time: Dec 16 04:08:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 14 02:24:06 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.387	152	11695	20.000	ng	0.01
21) Naphthalene-d8	11.219	136	43375	20.000	ng	# 0.01
39) Acenaphthene-d10	14.991	164	33468	20.000	ng	0.01
64) Phenanthrene-d10	17.735	188	75738	20.000	ng	0.02
76) Chrysene-d12	22.048	240	79195	20.000	ng	0.02
86) Perylene-d12	25.561	264	60033	20.000	ng	# 0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.896	112	9330	18.709	ng	0.01
7) Phenol-d6	7.506	99	13521	21.081	ng	0.01
23) Nitrobenzene-d5	9.550	82	9176	14.784	ng	0.00
42) 2,4,6-Tribromophenol	16.466	330	11663	14.452	ng	0.01
45) 2-Fluorobiphenyl	13.616	172	31745	13.303	ng	0.00
79) Terphenyl-d14	20.303	244	53167	11.865	ng	0.01
Target Compounds						Qvalue
37) 2-Methylnaphthalene	12.847	142	27847	13.772	ng	99
38) 1-Methylnaphthalene	13.058	142	23947	12.157	ng	97
46) 1,1'-Biphenyl	13.828	154	9943	4.383	ng	98
52) Acenaphthene	15.056	154	39580	20.038	ng	95
55) Dibenzofuran	15.385	168	16239	4.889	ng	# 94
58) Fluorene	16.031	166	55849	20.569	ng	96
71) Phenanthrene	17.788	178	1109406	279.513	ng	98
72) Anthracene	17.864	178	158971	40.485	ng	99
75) Fluoranthene	19.768	202	550565	99.060	ng	98
78) Pyrene	20.132	202	829505	173.097	ng	99
81) Benzo(a)anthracene	22.024	228	341770	61.328	ng	99
83) Chrysene	22.100	228	343226	65.716	ng	98
87) Indeno(1,2,3-cd)pyrene	29.592	276	32957m	6.827	ng	
88) Benzo(b)fluoranthene	24.445	252	139522	35.911	ng	# 94
89) Benzo(k)fluoranthene	24.498	252	95848m	24.726	ng	
90) Benzo(a)pyrene	25.391	252	120785	36.778	ng	97
91) Dibenzo(a,h)anthracene	29.645	278	12543m	3.104	ng	
92) Benzo(g,h,i)perylene	30.872	276	25016m	6.403	ng	

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

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